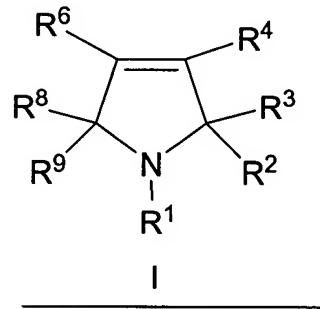
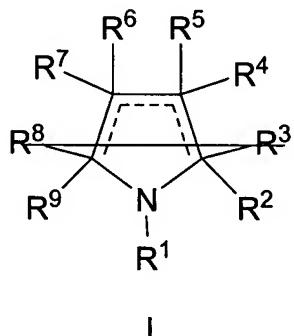


In the claims:

1. (Currently amended)

A compound of Formula I:



or a pharmaceutically acceptable salt or stereoisomer thereof, wherein

a is 0 or 1;

b is 0 or 1;

m is 0, 1, or 2;

n is 0 or 1;

r is 0 or 1;

s is 0 or 1;

u is 2, 3, 4 or 5;

~~a dashed line represents an optional double bond, provided that one and only one double bond is present in the ring;~~

R<sup>1</sup> is selected from:

- 1)  $(C_1-C_6\text{-alkylene})_n(C=X)C_1-C_{10}\text{ alkyl},$
- 2)  $(C_1-C_6\text{-alkylene})_n(C=X)\text{aryl},$
- 3)  $(C_1-C_6\text{-alkylene})_n(C=X)C_2-C_{10}\text{ alkenyl},$
- 4)  $(C_1-C_6\text{-alkylene})_n(C=X)C_2-C_{10}\text{ alkynyl},$
- 5)  $(C_1-C_6\text{-alkylene})_n(C=X)C_3-C_8\text{ cycloalkyl},$
- 6)  $(C_1-C_6\text{-alkylene})_n(C=X)\text{heterocyclyl},$

- 7)  $(C_1\text{-}C_6\text{-alkylene})_n(C=X)NR^cR^{c'},$
- 8)  $(C_1\text{-}C_6\text{-alkylene})_nSO_2NR^eR^{e'},$
- 9)  $(C_1\text{-}C_6\text{-alkylene})_nSO_2C_1\text{-}C_{10}\text{-alkyl},$
- 10)  $(C_1\text{-}C_6\text{-alkylene})_nSO_2C_2\text{-}C_{10}\text{-alkenyl},$
- 11)  $(C_1\text{-}C_6\text{-alkylene})_nSO_2C_2\text{-}C_{10}\text{-alkynyl},$
- 12)  $(C_1\text{-}C_6\text{-alkylene})_nSO_2\text{-aryl},$
- 13)  $(C_1\text{-}C_6\text{-alkylene})_nSO_2\text{-heterocyclyl},$
- 14)  $(C_1\text{-}C_6\text{-alkylene})_nSO_2\text{-}C_3\text{-}C_8\text{-cycloalkyl},$
- 15)  $(C_1\text{-}C_6\text{-alkylene})_nP(=O)R^dR^{d'},$
- 16) aryl;
- 17) heterocyclyl; and
- 18)  $C_1\text{-}C_{10}\text{ alkyl};$

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, alkylene, heteroaryl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>2</sup> and R<sup>6</sup> are independently selected from:

- 1) aryl,
- 2)  $C_1\text{-}C_6\text{-aralkyl},$
- 3)  $C_3\text{-}C_8\text{-cycloalkyl},$  and
- 4) heterocyclyl,

said aryl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>3</sup>, R<sup>4</sup>, R<sup>5</sup>, R<sup>7</sup>, R<sup>8</sup>, and R<sup>9</sup> are independently selected from:

- 1) H, and
- 2)  $C_1\text{-}C_{10} \text{ alkyl},$
- 3) aryl,
- 4)  $C_2\text{-}C_{10} \text{ alkenyl},$
- 5)  $C_2\text{-}C_{10} \text{ alkynyl},$
- 7)  $C_1\text{-}C_6 \text{ perfluoroalkyl},$
- 8)  $C_1\text{-}C_6 \text{ aralkyl},$

9) ~~C<sub>3</sub>-C<sub>8</sub> cycloalkyl, and~~

10) ~~heterocyclyl,~~

~~said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>, or~~

~~R<sup>4</sup> and R<sup>5</sup>, or R<sup>8</sup> and R<sup>9</sup>, attached to the same carbon atom are combined to form -(CH<sub>2</sub>)<sub>n</sub> wherein one of the carbon atoms is optionally replaced by a moiety selected from O, S(O)<sub>m</sub>, N(R<sup>a</sup>)C(O), N(R<sup>b</sup>) and N(COR<sup>a</sup>);~~

R<sup>10</sup> is independently selected from:

- 1) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 2) (C=O)<sub>a</sub>O<sub>b</sub>aryl,
- 3) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 4) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 5) (C=O)<sub>a</sub>O<sub>b</sub> heterocyclyl,
- 6) CO<sub>2</sub>H,
- 7) halo,
- 8) CN,
- 9) OH,
- 10) O<sub>b</sub>C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,
- 11) O<sub>a</sub>(C=O)<sub>b</sub>NR<sup>12</sup>R<sup>13</sup>,
- 12) S(O)<sub>m</sub>R<sup>a</sup>,
- 13) S(O)<sub>2</sub>NR<sup>12</sup>R<sup>13</sup>,
- 14) ~~exo;~~
- 15) CHO, ~~and~~
- 16) ~~(N=O)R<sup>12</sup>R<sup>13</sup>, or~~
- 17) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>3</sub>-C<sub>8</sub> cycloalkyl,

~~said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one or more substituents selected from R<sup>11</sup>;~~

R<sup>11</sup> is selected from:

- 1) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>1</sub>-C<sub>10</sub>)alkyl,

- 2)  $O_r(C_1-C_3)$ perfluoroalkyl,
- 3)  $(C_0-C_6)$ alkylene-S(O)<sub>m</sub>R<sup>a</sup>,
- 4) oxo,
- 5) OH,
- 6) halo,
- 7) CN,
- 8)  $(C=O)_rO_s(C_2-C_{10})$ alkenyl,
- 9)  $(C=O)_rO_s(C_2-C_{10})$ alkynyl,
- 10)  $(C=O)_rO_s(C_3-C_6)$ cycloalkyl,
- 11)  $(C=O)_rO_s(C_0-C_6)$ alkylene-aryl,
- 12)  $(C=O)_rO_s(C_0-C_6)$ alkylene-heterocyclyl,
- 13)  $(C=O)_rO_s(C_0-C_6)$ alkylene-N(R<sup>b</sup>)<sub>2</sub>,
- 14) C(O)R<sup>a</sup>,
- 15)  $(C_0-C_6)$ alkylene-CO<sub>2</sub>R<sup>a</sup>,
- 16) C(O)H,
- 17)  $(C_0-C_6)$ alkylene-CO<sub>2</sub>H,
- 18) C(O)N(R<sup>b</sup>)<sub>2</sub>,
- 19) S(O)<sub>m</sub>R<sup>a</sup>, and
- 20) S(O)<sub>2</sub>N(R<sup>b</sup>)<sub>2</sub>

said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, alkylene and heterocyclyl is optionally substituted with up to three substituents selected from R<sup>b</sup>, OH,  $(C_1-C_6)$ alkoxy, halogen, CO<sub>2</sub>H, CN, O(C=O)C<sub>1</sub>-C<sub>6</sub> alkyl, oxo, and N(R<sup>b</sup>)<sub>2</sub>;

R<sup>12</sup> and R<sup>13</sup> are independently selected from:

- 1) H,
- 2)  $(C=O)O_bC_1-C_{10}$  alkyl,
- 3)  $(C=O)O_bC_3-C_8$  cycloalkyl,
- 4)  $(C=O)O_b$ aryl,
- 5)  $(C=O)O_b$ heterocyclyl,
- 6) C<sub>1</sub>-C<sub>10</sub> alkyl,
- 7) aryl,
- 8) C<sub>2</sub>-C<sub>10</sub> alkenyl,

- 9) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 10) heterocyclyl,
- 11) C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
- 12) SO<sub>2</sub>R<sup>a</sup>, and
- 13) (C=O)NR<sup>b</sup><sub>2</sub>,

said alkyl, cycloalkyl, aryl, heterocyclyl, alkenyl, and alkynyl is optionally substituted with one or more substituents selected from R<sup>11</sup>, or

R<sup>12</sup> and R<sup>13</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 3-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one or more substituents selected from R<sup>11</sup>;

R<sup>14</sup> is independently selected from:

- 1) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 2) (C=O)<sub>a</sub>O<sub>b</sub>aryl,
- 3) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 4) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 5) (C=O)<sub>a</sub>O<sub>b</sub> heterocyclyl,
- 6) CO<sub>2</sub>H,
- 7) halo,
- 8) CN,
- 9) OH,
- 10) O<sub>b</sub>C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,
- 11) O<sub>a</sub>(C=O)<sub>b</sub>NR<sup>12</sup>R<sup>13</sup>,
- 12) S(O)<sub>m</sub>R<sup>a</sup>,
- 13) S(O)<sub>2</sub>NR<sup>12</sup>R<sup>13</sup>,
- 14) oxo,
- 15) CHO,
- 16) (N=O)R<sup>12</sup>R<sup>13</sup>, or
- 17) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>3</sub>-C<sub>8</sub> cycloalkyl,

said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one or more substituents selected from R<sup>11</sup>;

R<sup>a</sup> is (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, aryl, or heterocyclyl, optionally substituted with one to three substituents selected from R<sup>14</sup>;

R<sup>b</sup> is H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, (C=O)OC<sub>1</sub>-C<sub>6</sub> alkyl, (C=O)C<sub>1</sub>-C<sub>6</sub> alkyl or S(O)<sub>2</sub>R<sup>a</sup>, optionally substituted with one to three substituents selected from R<sup>14</sup>;

R<sup>c</sup> and R<sup>c'</sup> are independently selected from: H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl and (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, optionally substituted with one, two or three substituents selected from R<sup>10</sup>, or R<sup>c</sup> and R<sup>c'</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 3-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>;

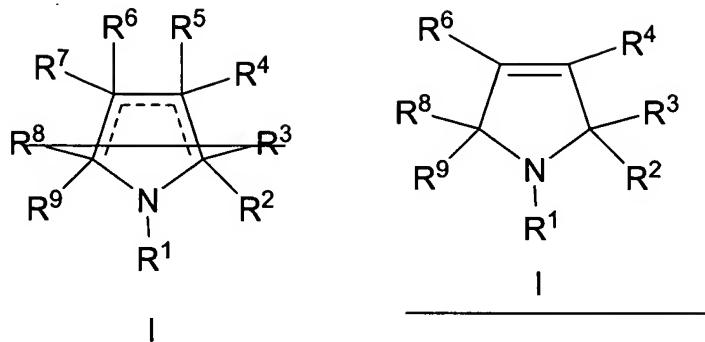
R<sup>d</sup> and R<sup>d'</sup> are independently selected from: (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>1</sub>-C<sub>6</sub>)alkoxy and NR<sup>b</sup><sub>2</sub>, or

R<sup>d</sup> and R<sup>d'</sup> can be taken together with the phosphorous to which they are attached to form a monocyclic heterocycle with 5-7 members in the ring and optionally containing, in addition to the phosphorous, one or two additional heteroatoms selected from NR<sup>e</sup>, O and S, said monocyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>;

R<sup>e</sup> is selected from: H and (C<sub>1</sub>-C<sub>6</sub>)alkyl; and

X is selected from O, NR<sup>e</sup> and S.

2. (Currently amended) The compound according to Claim 1 of the Formula I:



or a pharmaceutically acceptable salt or stereoisomer thereof, wherein

a is 0 or 1;  
b is 0 or 1;  
m is 0, 1, or 2;  
n is 0 or 1;  
r is 0 or 1;  
s is 0 or 1;  
u is 2, 3, 4 or 5;

~~a dashed line represents an optional double bond, provided that one and only one double bond is present in the ring;~~

R<sup>1</sup> is selected from:

- 1) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>1</sub>-C<sub>10</sub> alkyl,
- 2) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)aryl,
- 3) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 4) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 5) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
- 6) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)heterocyclyl,
- 7) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)NR<sup>c</sup>RC',
- 8) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>NR<sup>e</sup>RE',
- 9) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 10) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>C<sub>2</sub>-C<sub>10</sub> alkenyl,

- 11)  $(C_1-C_6\text{-alkylene})_nSO_2C_2-C_{10}\text{-alkynyl}$ ,
- 12)  $(C_1-C_6\text{-alkylene})_nSO_2\text{-aryl}$ ,
- 13)  $(C_1-C_6\text{-alkylene})_nSO_2\text{-heterocyclyl}$ ,
- 14)  $(C_1-C_6\text{-alkylene})_nSO_2-C_3-C_8\text{-cycloalkyl}$ ,
- 15)  $(C_1-C_6\text{-alkylene})_nP(=O)R^dR^d'$ ,
- 16) aryl;
- 17) heterocyclyl; and
- 18)  $C_1-C_{10}\text{ alkyl}$ ;

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, alkylene, heteroaryl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>2</sup> and R<sup>6</sup> are independently selected from:

- 1) aryl,
- 2)  $C_1-C_6\text{-aralkyl}$ ,
- 3)  $C_3-C_8\text{-cycloalkyl}$ , and
- 4) heterocyclyl,

said aryl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>3</sup>, R<sup>4</sup>, R<sup>5</sup>, R<sup>7</sup>, R<sup>8</sup>, and R<sup>9</sup> are independently selected from:

- 1) H, and
- 2)  $C_1-C_{10}\text{ alkyl}$ ,
- 3) aryl,
- 4)  $C_2-C_{10}\text{-alkenyl}$ ,
- 5)  $C_2-C_{10}\text{-alkynyl}$ ,
- 6)  $C_1-C_6\text{-perfluoroalkyl}$ ,
- 7)  $C_1-C_6\text{-aralkyl}$ ,
- 8)  $C_3-C_8\text{-cycloalkyl}$ , and
- 9) heterocyclyl,

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>, or

~~R<sup>4</sup> and R<sup>5</sup>, or R<sup>8</sup> and R<sup>9</sup>, attached to the same carbon atom are combined to form -(CH<sub>2</sub>)<sub>n</sub>- wherein one of the carbon atoms is optionally replaced by a moiety selected from O, S(O)<sub>m</sub>, N(R<sup>a</sup>)C(O), N(R<sup>b</sup>) and N(COR<sup>a</sup>);~~

R<sup>10</sup> is independently selected from:

- 1) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 2) (C=O)<sub>a</sub>O<sub>b</sub>aryl,
- 3) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 4) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 5) (C=O)<sub>a</sub>O<sub>b</sub> heterocyclyl,
- 6) CO<sub>2</sub>H,
- 7) halo,
- 8) CN,
- 9) OH,
- 10) O<sub>b</sub>C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,
- 11) O<sub>a</sub>(C=O)<sub>b</sub>NR<sup>12</sup>R<sup>13</sup>,
- 12) S(O)<sub>m</sub>R<sup>a</sup>,
- 13) S(O)<sub>2</sub>NR<sup>12</sup>R<sup>13</sup>,
- 14) ~~oxo,~~
- 15) CHO, and
- 16) ~~(N=O)R<sup>12</sup>R<sup>13</sup>, or~~
- 17) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>3</sub>-C<sub>8</sub> cycloalkyl,

said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one or more substituents selected from R<sup>11</sup>;

R<sup>11</sup> is selected from:

- 1) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>1</sub>-C<sub>10</sub>)alkyl,
- 2) O<sub>r</sub>(C<sub>1</sub>-C<sub>3</sub>)perfluoroalkyl,
- 3) (C<sub>0</sub>-C<sub>6</sub>)alkylene-S(O)<sub>m</sub>R<sup>a</sup>,
- 4) oxo,
- 5) OH,
- 6) halo,

- 7) CN,
- 8) (C=O)<sub>r</sub>Os(C<sub>2</sub>-C<sub>10</sub>)alkenyl,
- 9) (C=O)<sub>r</sub>Os(C<sub>2</sub>-C<sub>10</sub>)alkynyl,
- 10) (C=O)<sub>r</sub>Os(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
- 11) (C=O)<sub>r</sub>Os(C<sub>0</sub>-C<sub>6</sub>)alkylene-aryl,
- 12) (C=O)<sub>r</sub>Os(C<sub>0</sub>-C<sub>6</sub>)alkylene-heterocyclyl,
- 13) (C=O)<sub>r</sub>Os(C<sub>0</sub>-C<sub>6</sub>)alkylene-N(R<sup>b</sup>)<sub>2</sub>,
- 14) C(O)R<sup>a</sup>,
- 15) (C<sub>0</sub>-C<sub>6</sub>)alkylene-CO<sub>2</sub>R<sup>a</sup>,
- 16) C(O)H,
- 17) (C<sub>0</sub>-C<sub>6</sub>)alkylene-CO<sub>2</sub>H,
- 18) C(O)N(R<sup>b</sup>)<sub>2</sub>,
- 19) S(O)<sub>m</sub>R<sup>a</sup>, and
- 20) S(O)<sub>2</sub>N(R<sup>b</sup>)<sub>2</sub>

said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, alkylene and heterocyclyl is optionally substituted with up to three substituents selected from R<sup>b</sup>, OH, (C<sub>1</sub>-C<sub>6</sub>)alkoxy, halogen, CO<sub>2</sub>H, CN, O(C=O)C<sub>1</sub>-C<sub>6</sub> alkyl, oxo, and N(R<sup>b</sup>)<sub>2</sub>;

R<sup>12</sup> and R<sup>13</sup> are independently selected from:

- 1) H,
- 2) (C=O)O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 3) (C=O)O<sub>b</sub>C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
- 4) (C=O)Obaryl,
- 5) (C=O)Obheterocyclyl,
- 6) C<sub>1</sub>-C<sub>10</sub> alkyl,
- 7) aryl,
- 8) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 9) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 10) heterocyclyl,
- 11) C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
- 12) SO<sub>2</sub>R<sup>a</sup>, and
- 13) (C=O)NR<sup>b</sup><sub>2</sub>,

said alkyl, cycloalkyl, aryl, heterocyl, alkenyl, and alkynyl is optionally substituted with one or more substituents selected from R<sup>11</sup>, or

R<sup>12</sup> and R<sup>13</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one or more substituents selected from R<sup>11</sup>;

R<sup>a</sup> is (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, aryl, or heterocyclyl;

R<sup>b</sup> is H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, (C=O)OC<sub>1</sub>-C<sub>6</sub> alkyl, (C=O)C<sub>1</sub>-C<sub>6</sub> alkyl or S(O)<sub>2</sub>R<sup>a</sup>;

R<sup>c</sup> and R<sup>c'</sup> are independently selected from: H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl and (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, optionally substituted with one, two or three substituents selected from R<sup>10</sup>, or

R<sup>c</sup> and R<sup>c'</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>; and

~~R<sup>d</sup> and R<sup>d'</sup> are independently selected from: (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>1</sub>-C<sub>6</sub>)alkoxy and NR<sup>b</sup><sub>2</sub>, or~~

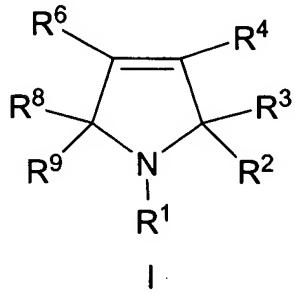
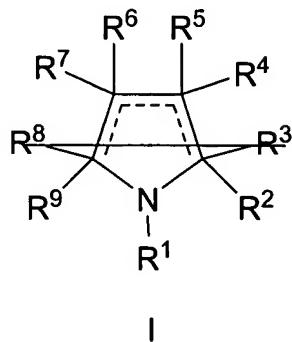
~~R<sup>d</sup> and R<sup>d'</sup> can be taken together with the phosphorous to which they are attached to form a monocyclic heterocycle with 5-7 members the ring and optionally containing, in addition to the phosphorous, one or two additional heteroatoms selected from NR<sup>e</sup>, O and S, said monocyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>;~~

R<sup>e</sup> is selected from: H and (C<sub>1</sub>-C<sub>6</sub>)alkyl; and

X is selected from O, NR<sup>a</sup> and S.

3. (Currently amended)

The compound according to Claim 2 of Formula I:



or a pharmaceutically acceptable salt or stereoisomer thereof, wherein

a is 0 or 1;  
b is 0 or 1;  
m is 0, 1, or 2;  
n is 0 or 1;  
r is 0 or 1;  
s is 0 or 1;  
~~u is 2, 3, 4 or 5;~~

~~a dashed line represents an optional double bond, provided that one and only one double bond is present in the ring;~~

R<sup>1</sup> is selected from:

- 1) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>1</sub>-C<sub>10</sub>-alkyl,
- 2) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)aryl,
- 3) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>2</sub>-C<sub>10</sub>-alkenyl,
- 4) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>2</sub>-C<sub>10</sub>-alkynyl,
- 5) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>3</sub>-C<sub>8</sub>-cycloalkyl,
- 6) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)heterocyclyl,

- 7)  $(C_1\text{-}C_6\text{-alkylene})_n(C=X)NR^cR^{c'},$
- 8)  $(C_1\text{-}C_6\text{-alkylene})_nSO_2NReR^{e'},$
- 9)  $(C_1\text{-}C_6\text{-alkylene})_nSO_2C_1\text{-}C_{10}\text{ alkyl},$
- 10)  $(C_1\text{-}C_6\text{-alkylene})_nSO_2C_2\text{-}C_{10}\text{ alkenyl},$
- 11)  $(C_1\text{-}C_6\text{-alkylene})_nSO_2C_2\text{-}C_{10}\text{ alkynyl},$
- 12)  $(C_1\text{-}C_6\text{-alkylene})_nSO_2\text{-aryl},$
- 13)  $(C_1\text{-}C_6\text{-alkylene})_nSO_2\text{-heterocyclyl},$
- 14)  $(C_1\text{-}C_6\text{-alkylene})_nSO_2\text{-}C_3\text{-}C_8\text{-cycloalkyl},$
- 15)  $(C_1\text{-}C_6\text{-alkylene})_nP(=O)R^dR^{d'},$
- 16) aryl;
- 17) heterocyclyl; and
- 18)  $C_1\text{-}C_{10}\text{ alkyl};$

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, alkylene, heteroaryl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>2</sup> and R<sup>6</sup> are independently selected from:

- 1) aryl,
- 2)  $C_1\text{-}C_6\text{-aralkyl},$
- 3)  $C_3\text{-}C_8\text{-cycloalkyl},$  and
- 4) heterocyclyl,

said aryl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>3</sup>, R<sup>4</sup>, R<sup>5</sup>, R<sup>7</sup>, R<sup>8</sup>, and R<sup>9</sup> are independently selected from:

- 1) H, and
- 2)  $C_1\text{-}C_{10}\text{ alkyl},$
- 3) aryl,
- 4)  $C_2\text{-}C_{10}\text{ alkenyl},$
- 5)  $C_2\text{-}C_{10}\text{ alkynyl},$
- 6)  $C_1\text{-}C_6\text{-perfluoroalkyl},$
- 7)  $C_1\text{-}C_6\text{-aralkyl},$
- 8)  $C_3\text{-}C_8\text{-cycloalkyl},$  and

9) heterocyclyl,

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>, or

R<sup>4</sup> and R<sup>5</sup>, or R<sup>8</sup> and R<sup>9</sup>, attached to the same carbon atom are combined to form -(CH<sub>2</sub>)<sub>n</sub> wherein one of the carbon atoms is optionally replaced by a moiety selected from O, S(O)<sub>m</sub>, N(R<sup>a</sup>)C(O), N(R<sup>b</sup>) and N(COR<sup>a</sup>);

R<sup>10</sup> is independently selected from:

- 1) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 2) (C=O)<sub>a</sub>O<sub>b</sub>aryl,
- 3) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 4) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 5) (C=O)<sub>a</sub>O<sub>b</sub> heterocyclyl,
- 6) CO<sub>2</sub>H,
- 7) halo,
- 8) CN,
- 9) OH,
- 10) O<sub>b</sub>C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,
- 11) O<sub>a</sub>(C=O)<sub>b</sub>NR<sup>12</sup>R<sup>13</sup>,
- 12) S(O)<sub>m</sub>R<sup>a</sup>,
- 13) S(O)<sub>2</sub>NR<sup>12</sup>R<sup>13</sup>,
- 14) -exo,
- 15) CHO, and
- 16) (N=O)R<sup>12</sup>R<sup>13</sup>, or
- 17) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>3</sub>-C<sub>8</sub> cycloalkyl,

said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one or more substituents selected from R<sup>11</sup>;

R<sup>11</sup> is selected from:

- 1) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>1</sub>-C<sub>10</sub>)alkyl,
- 2) O<sub>r</sub>(C<sub>1</sub>-C<sub>3</sub>)perfluoroalkyl,
- 3) (C<sub>0</sub>-C<sub>6</sub>)alkylene-S(O)<sub>m</sub>R<sup>a</sup>,

- 4) oxo,
- 5) OH,
- 6) halo,
- 7) CN,
- 8) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>2</sub>-C<sub>10</sub>)alkenyl,
- 9) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>2</sub>-C<sub>10</sub>)alkynyl,
- 10) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
- 11) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-aryl,
- 12) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-heterocyclyl,
- 13) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-N(R<sup>b</sup>)<sub>2</sub>,
- 14) C(O)R<sup>a</sup>,
- 15) (C<sub>0</sub>-C<sub>6</sub>)alkylene-CO<sub>2</sub>R<sup>a</sup>,
- 16) C(O)H,
- 17) (C<sub>0</sub>-C<sub>6</sub>)alkylene-CO<sub>2</sub>H,
- 18) C(O)N(R<sup>b</sup>)<sub>2</sub>,
- 19) S(O)<sub>m</sub>R<sup>a</sup>, and
- 20) S(O)<sub>2</sub>N(R<sup>b</sup>)<sub>2</sub>,

said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, alkylene and heterocyclyl is optionally substituted with up to three substituents selected from R<sup>b</sup>, OH, (C<sub>1</sub>-C<sub>6</sub>)alkoxy, halogen, CO<sub>2</sub>H, CN, O(C=O)C<sub>1</sub>-C<sub>6</sub> alkyl, oxo, and N(R<sup>b</sup>)<sub>2</sub>;

R<sup>12</sup> and R<sup>13</sup> are independently selected from:

- 1) H,
- 2) (C=O)O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 3) (C=O)O<sub>b</sub>C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
- 4) (C=O)O<sub>b</sub>aryl,
- 5) (C=O)O<sub>b</sub>heterocyclyl,
- 6) C<sub>1</sub>-C<sub>10</sub> alkyl,
- 7) aryl,
- 8) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 9) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 10) heterocyclyl,

- 11) C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
- 12) SO<sub>2</sub>R<sup>a</sup>, and
- 13) (C=O)NR<sup>b</sup><sub>2</sub>,

said alkyl, cycloalkyl, aryl, heterocyl, alkenyl, and alkynyl is optionally substituted with one or more substituents selected from R<sup>11</sup>, or

R<sup>12</sup> and R<sup>13</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one or more substituents selected from R<sup>11</sup>;

R<sup>a</sup> is (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, aryl, or heterocyclyl;

R<sup>b</sup> is H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, (C=O)OC<sub>1</sub>-C<sub>6</sub> alkyl, (C=O)C<sub>1</sub>-C<sub>6</sub> alkyl or S(O)<sub>2</sub>R<sup>a</sup>;

R<sup>c</sup> and R<sup>c'</sup> are independently selected from: H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl and (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl or

R<sup>c</sup> and R<sup>c'</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>;

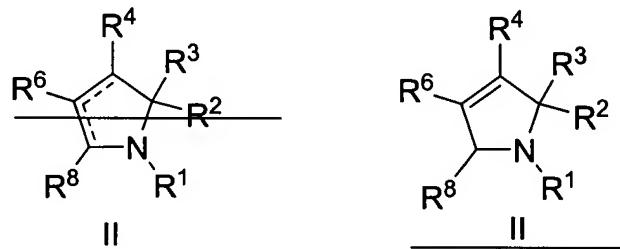
~~R<sup>d</sup> and R<sup>d'</sup> are independently selected from: (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>1</sub>-C<sub>6</sub>)alkoxy and NR<sup>b</sup><sub>2</sub>, or~~

~~R<sup>d</sup> and R<sup>d'</sup> can be taken together with the phosphorous to which they are attached to form a monocyclic heterocycle with 5-7 members the ring and optionally containing, in addition to the phosphorous, one or two additional heteroatoms selected from N<sup>e</sup>, O and S, said monocyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>;~~

R<sup>e</sup> is selected from: H and (C<sub>1</sub>-C<sub>6</sub>)alkyl; and

X is selected from O, NR<sup>e</sup> and S.

4. (Currently amended) The compound according to Claim 2 of the Formula II,



or a pharmaceutically acceptable salt or stereoisomer thereof, wherein  
wherein:

a is 0 or 1;  
b is 0 or 1;  
m is 0, 1, or 2;  
r is 0 or 1;  
s is 0 or 1;

~~a dashed line represents an optional double bond, provided that one and only one double bond is present in the ring;~~

R<sup>1</sup> is selected from:

- 1) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=O)C<sub>1</sub>-C<sub>10</sub>-alkyl,
- 2) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=O)aryl,
- 3) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=O)C<sub>2</sub>-C<sub>10</sub>-alkenyl,
- 4) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=O)C<sub>2</sub>-C<sub>10</sub>-alkynyl,
- 5) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=O)C<sub>3</sub>-C<sub>8</sub>-cycloalkyl,
- 6) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=O)heterocyclyl,

- 7)  $(C_1\text{-}C_6\text{-alkylene})_n(C=O)NR^cR^c'$ ,
- 8)  $(C_1\text{-}C_6\text{-alkylene})_nSO_2NR^eR^e'$ ,
- 9)  $(C_1\text{-}C_6\text{-alkylene})_nSO_2C_1\text{-}C_{10}\text{-alkyl}$ ,
- 10)  $(C_1\text{-}C_6\text{-alkylene})_nSO_2\text{-aryl}$ ,
- 11)  $(C_1\text{-}C_6\text{-alkylene})_nSO_2\text{-heterocyclyl}$ ,
- 12)  $(C_1\text{-}C_6\text{-alkylene})_nSO_2\text{-}C_3\text{-}C_8\text{-cycloalkyl}$ ,
- 13)  $(C_1\text{-}C_6\text{-alkylene})_nP(=O)R^dR^d'$ ,
- 14)  $\text{aryl}$ ,
- 15) heterocyclyl, and
- 16)  $C_1\text{-}C_{10}\text{-alkyl}$ ;

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, alkylene, heteroaryl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>2</sup> and R<sup>6</sup> are independently selected from:

- 1)  $\text{aryl}$ ,
- 2)  $C_1\text{-}C_6\text{-aralkyl}$ ,
- 3)  $C_3\text{-}C_8\text{-cycloalkyl}$ , and
- 4) heterocyclyl,

said aryl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>3</sup>, R<sup>4</sup> and R<sup>8</sup> are independently selected from:

- 1) H, and
- 2)  $C_1\text{-}C_{10}\text{ alkyl}$ ,
- 3)  $\text{aryl}$ ,
- 4)  $C_2\text{-}C_{10}\text{-alkenyl}$ ,
- 5)  $C_2\text{-}C_{10}\text{-alkynyl}$ ,
- 6)  $C_1\text{-}C_6\text{-perfluoroalkyl}$ ,
- 7)  $C_1\text{-}C_6\text{-aralkyl}$ ,
- 8)  $C_3\text{-}C_8\text{-cycloalkyl}$ , and
- 9) heterocyclyl,

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>10</sup> is independently selected from:

- 1) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 2) (C=O)<sub>a</sub>O<sub>b</sub>aryl,
- 3) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 4) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 5) (C=O)<sub>a</sub>O<sub>b</sub> heterocyclyl,
- 6) CO<sub>2</sub>H,
- 7) halo,
- 8) CN,
- 9) OH,
- 10) O<sub>b</sub>C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,
- 11) O<sub>a</sub>(C=O)<sub>b</sub>NR<sub>12</sub>R<sub>13</sub>,
- 12) S(O)<sub>m</sub>R<sup>a</sup>,
- 13) S(O)<sub>2</sub>NR<sub>12</sub>R<sub>13</sub>,
- 14) ~~oxo,~~
- 15) CHO, and
- 16) ~~(N=O)R<sup>12</sup>R<sup>13</sup>, or~~
- 17) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>3</sub>-C<sub>8</sub> cycloalkyl,

said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one, two or three substituents selected from R<sup>11</sup>;

R<sup>11</sup> is selected from:

- 1) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>1</sub>-C<sub>10</sub>)alkyl,
- 2) O<sub>r</sub>(C<sub>1</sub>-C<sub>3</sub>)perfluoroalkyl,
- 3) oxo,
- 4) OH,
- 5) halo,
- 6) CN,
- 7) (C<sub>2</sub>-C<sub>10</sub>)alkenyl,

- 8) (C<sub>2</sub>-C<sub>10</sub>)alkynyl,
- 9) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
- 10) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-aryl,
- 11) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-heterocyclyl,
- 12) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-N(R<sup>b</sup>)<sub>2</sub>,
- 13) C(O)R<sup>a</sup>,
- 14) (C<sub>0</sub>-C<sub>6</sub>)alkylene-CO<sub>2</sub>R<sup>a</sup>,
- 15) C(O)H,
- 16) (C<sub>0</sub>-C<sub>6</sub>)alkylene-CO<sub>2</sub>H,
- 17) C(O)N(R<sup>b</sup>)<sub>2</sub>,
- 18) S(O)<sub>m</sub>R<sup>a</sup>, and
- 19) S(O)<sub>2</sub>N(R<sup>b</sup>)<sub>2</sub>;

said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, alkylene and heterocyclyl is optionally substituted with up to three substituents selected from R<sup>b</sup>, OH, (C<sub>1</sub>-C<sub>6</sub>)alkoxy, halogen, CO<sub>2</sub>H, CN, O(C=O)C<sub>1</sub>-C<sub>6</sub> alkyl, oxo, and N(R<sup>b</sup>)<sub>2</sub>;

R<sup>12</sup> and R<sup>13</sup> are independently selected from:

- 1) H,
- 2) (C=O)O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 3) (C=O)O<sub>b</sub>C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
- 4) (C=O)O<sub>b</sub>aryl,
- 5) (C=O)O<sub>b</sub>heterocyclyl,
- 6) C<sub>1</sub>-C<sub>10</sub> alkyl,
- 7) aryl,
- 8) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 9) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 10) heterocyclyl,
- 11) C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
- 12) SO<sub>2</sub>R<sup>a</sup>, and
- 13) (C=O)NR<sup>b</sup><sub>2</sub>,

said alkyl, cycloalkyl, aryl, heterocylyl, alkenyl, and alkynyl is optionally substituted with one, two or three substituents selected from R<sup>11</sup>, or

R<sup>12</sup> and R<sup>13</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>;

R<sup>a</sup> is (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, aryl, or heterocyclyl;

R<sup>b</sup> is H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, (C=O)OC<sub>1</sub>-C<sub>6</sub> alkyl, (C=O)C<sub>1</sub>-C<sub>6</sub> alkyl or S(O)<sub>2</sub>R<sup>a</sup>;

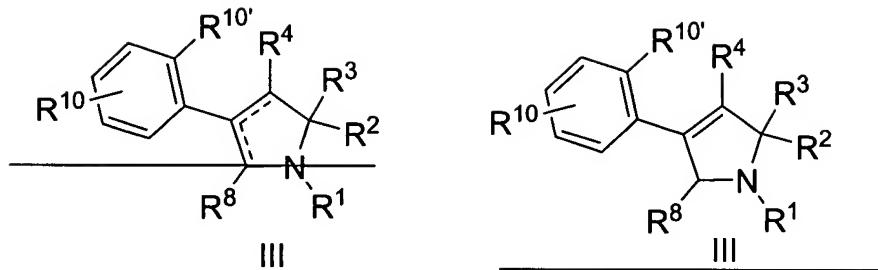
R<sup>c</sup> and R<sup>c'</sup> are independently selected from: H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl and (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl; or R<sup>c</sup> and R<sup>c'</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>; .

R<sup>d</sup> and R<sup>d'</sup> are independently selected from: (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>1</sub>-C<sub>6</sub>)alkoxy and NR<sup>b</sup><sub>2</sub>, or

~~R<sup>d</sup> and R<sup>d'</sup> can be taken together with the phosphorous to which they are attached to form a monocyclic heterocycle with 5-7 members the ring and optionally containing, in addition to the phosphorous, one or two additional heteroatoms selected from NR<sup>e</sup>, O and S, said monocyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>; and~~

R<sup>e</sup> is selected from: H and (C<sub>1</sub>-C<sub>6</sub>)alkyl.

5. (Currently amended) A compound of the Formula III,



or a pharmaceutically acceptable salt or stereoisomer thereof, wherein

a is 0 or 1;  
b is 0 or 1;  
m is 0, 1, or 2;  
r is 0 or 1;  
s is 0 or 1;

~~a dashed line represents an optional double bond, provided that one and only one double bond is present in the ring;~~

R<sup>1</sup> is selected from:

- 1)  $(C=O)C_1-C_{10}$ -alkyl,
- 2)  $(C=O)$ aryl,
- 3)  $(C=O)C_2-C_{10}$ -alkenyl,
- 4)  $(C=O)C_2-C_{10}$ -alkynyl,
- 5)  $(C=O)C_3-C_8$ -cycloalkyl,
- 6)  $(C=O)$ heterocyclyl,
- 7)  $(C=O)NR^cR^{c'}$ ,
- 8)  $SO_2NR^eR^{e'}$ ,
- 9)  $SO_2C_1-C_{10}$ -alkyl,
- 10)  $SO_2$ -aryl,
- 11)  $SO_2$ -heterocyclyl,
- 12)  $SO_2C_3-C_8$ -cycloalkyl, and
- 13)  $P(=O)R^dR^{d'}$ ,

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, heteroaryl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>2</sup> is selected from:

- 1) aryl,
- 2) C<sub>1</sub>-C<sub>6</sub> aralkyl,
- 3) C<sub>3</sub>-C<sub>8</sub> cycloalkyl, and
- 4) heterocyclyl,

said aryl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>3</sup>, R<sup>4</sup> and R<sup>8</sup> are independently selected from:

- 1) H, and
- 2) C<sub>1</sub>-C<sub>10</sub> alkyl,
- 3) aryl,
- 4) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 5) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 6) C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,
- 7) C<sub>1</sub>-C<sub>6</sub> aralkyl,
- 8) C<sub>3</sub>-C<sub>8</sub> cycloalkyl, and
- 9) heterocyclyl,

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>10</sup> is independently selected from:

- 1) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 2) (C=O)<sub>a</sub>O<sub>b</sub>aryl,
- 3) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 4) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 5) (C=O)<sub>a</sub>O<sub>b</sub> heterocyclyl,
- 6) CO<sub>2</sub>H,
- 7) halo,
- 8) CN,

- 9) OH,
- 10) O<sub>b</sub>C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,
- 11) O<sub>a</sub>(C=O)<sub>b</sub>NR<sup>12</sup>R<sup>13</sup>,
- 12) S(O)<sub>m</sub>R<sup>a</sup>,
- 13) S(O)<sub>2</sub>NR<sup>12</sup>R<sup>13</sup>,
- 14) ~~oxo;~~
- 15) CHO, and
- 16) ~~(N=O)R<sup>12</sup>R<sup>13</sup>, or~~
- 17) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>3</sub>-C<sub>8</sub> cycloalkyl,

said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one, two or three substituents selected from R<sup>11</sup>;

R<sup>10'</sup> is halogen;

R<sup>11</sup> is selected from:

- 1) (C=O)<sub>r</sub>Os(C<sub>1</sub>-C<sub>10</sub>)alkyl,
- 2) O<sub>r</sub>(C<sub>1</sub>-C<sub>3</sub>)perfluoroalkyl,
- 3) oxo,
- 4) OH,
- 5) halo,
- 6) CN,
- 7) (C<sub>2</sub>-C<sub>10</sub>)alkenyl,
- 8) (C<sub>2</sub>-C<sub>10</sub>)alkynyl,
- 9) (C=O)<sub>r</sub>Os(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
- 10) (C=O)<sub>r</sub>Os(C<sub>0</sub>-C<sub>6</sub>)alkylene-aryl,
- 11) (C=O)<sub>r</sub>Os(C<sub>0</sub>-C<sub>6</sub>)alkylene-heterocyclyl,
- 12) (C=O)<sub>r</sub>Os(C<sub>0</sub>-C<sub>6</sub>)alkylene-N(R<sup>b</sup>)<sub>2</sub>,
- 13) C(O)R<sup>a</sup>,
- 14) (C<sub>0</sub>-C<sub>6</sub>)alkylene-CO<sub>2</sub>R<sup>a</sup>,
- 15) C(O)H,
- 16) (C<sub>0</sub>-C<sub>6</sub>)alkylene-CO<sub>2</sub>H, and
- 17) C(O)N(R<sup>b</sup>)<sub>2</sub>,

- 18)  $S(O)_m R^a$ , and
- 19)  $S(O)2N(R^b)2$ ;

said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, and heterocyclyl is optionally substituted with up to three substituents selected from  $R^b$ , OH, ( $C_1-C_6$ )alkoxy, halogen,  $CO_2H$ , CN,  $O(C=O)C_1-C_6$  alkyl, oxo, and  $N(R^b)2$ ;

$R^{12}$  and  $R^{13}$  are independently selected from:

- 1) H,
- 2)  $(C=O)O_b C_1-C_{10}$  alkyl,
- 3)  $(C=O)O_b C_3-C_8$  cycloalkyl,
- 4)  $(C=O)O_b$  aryl,
- 5)  $(C=O)O_b$  heterocyclyl,
- 6)  $C_1-C_{10}$  alkyl,
- 7) aryl,
- 8)  $C_2-C_{10}$  alkenyl,
- 9)  $C_2-C_{10}$  alkynyl,
- 10) heterocyclyl,
- 11)  $C_3-C_8$  cycloalkyl,
- 12)  $SO_2R^a$ , and
- 13)  $(C=O)NR^b_2$ ,

said alkyl, cycloalkyl, aryl, heterocyclyl, alkenyl, and alkynyl is optionally substituted with one, two or three substituents selected from  $R^{11}$ , or

$R^{12}$  and  $R^{13}$  can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from  $R^{11}$ ;

$R^a$  is  $(C_1-C_6)$  alkyl,  $(C_3-C_6)$  cycloalkyl, aryl, or heterocyclyl;

R<sup>b</sup> is H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, (C=O)OC<sub>1</sub>-C<sub>6</sub> alkyl, (C=O)C<sub>1</sub>-C<sub>6</sub> alkyl or S(O)<sub>2</sub>R<sup>a</sup>;

R<sup>c</sup> and R<sup>c'</sup> are independently selected from: H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl and (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl; or

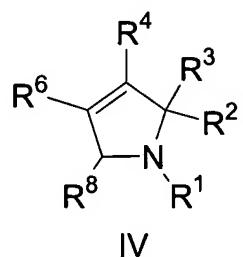
R<sup>c</sup> and R<sup>c'</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>;

R<sup>d</sup> and R<sup>d'</sup> are independently selected from: (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>1</sub>-C<sub>6</sub>)alkoxy and NR<sup>b</sup><sub>2</sub>, or

R<sup>d</sup> and R<sup>d'</sup> can be taken together with the phosphorous to which they are attached to form a monocyclic heterocycle with 5-7 members in the ring and optionally containing, in addition to the phosphorous, one or two additional heteroatoms selected from NR<sup>e</sup>, O and S, said monocyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>; and

R<sup>e</sup> is selected from: H and (C<sub>1</sub>-C<sub>6</sub>)alkyl.

6. (Currently amended) The compound according to Claim 4 of the Formula IV,



or a pharmaceutically acceptable salt or stereoisomer thereof, wherein

a is 0 or 1;  
b is 0 or 1;  
m is 0, 1, or 2;  
r is 0 or 1;  
s is 0 or 1;

R<sup>1</sup> is selected from:

- 1)  $(C=O)C_1-C_{10}$  alkyl,
- 2)  $(C=O)$ aryl,
- 3)  $(C=O)C_3-C_8$  cycloalkyl,
- 4)  $(C=O)$ heterocyclyl,
- 5)  $(C=O)NRcRC'$ ,
- 6)  $(C=S)NRcRC'$ ,
- 7)  $SO_2NRcRC'$ ,
- 8)  $SO_2C_1-C_{10}$  alkyl,
- 9)  $SO_2$  aryl, and
- 10)  $SO_2$  heterocyclyl,

said alkyl, aryl, cycloalkyl, and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>2</sup> is selected from:

- 1) aryl,
- 2)  $C_1-C_6$  aralkyl,
- 3)  $C_3-C_8$  cycloalkyl, and
- 4) heterocyclyl,

said aryl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>3</sup>, R<sup>4</sup> and R<sup>8</sup> are independently selected from:

- 1) H, and
- 2) C<sub>1</sub>-C<sub>10</sub> alkyl, and
- 3)  $C_1-C_6$  perfluoroalkyl,

~~said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;~~

R<sup>6</sup> is selected from:

- 1) aryl,
- 2) ~~C<sub>4</sub>-C<sub>6</sub> aralkyl,~~
- 3) ~~C<sub>3</sub>-C<sub>8</sub> cycloalkyl, and~~
- 4) ~~heterocyclyl,~~

~~said aryl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;~~

R<sup>10</sup> is independently selected from:

- 1) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 2) (C=O)<sub>a</sub>O<sub>b</sub>aryl,
- 3) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 4) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 5) (C=O)<sub>a</sub>O<sub>b</sub> heterocyclyl,
- 6) CO<sub>2</sub>H,
- 7) halo,
- 8) CN,
- 9) OH,
- 10) O<sub>b</sub>C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,
- 11) O<sub>a</sub>(C=O)<sub>b</sub>NR<sup>12</sup>R<sup>13</sup>,
- 12) S(O)<sub>m</sub>R<sup>a</sup>,
- 13) S(O)<sub>2</sub>NR<sup>12</sup>R<sup>13</sup>,
- 14) ~~exo-~~,
- 15) CHO, and
- 16) ~~(N=O)R<sup>12</sup>R<sup>13</sup>, or~~
- 17) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>3</sub>-C<sub>8</sub> cycloalkyl,

~~said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one, two or three substituents selected from R<sup>11</sup>;~~

R<sup>11</sup> is selected from:

- 1) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>1</sub>-C<sub>10</sub>)alkyl,
- 2) O<sub>r</sub>(C<sub>1</sub>-C<sub>3</sub>)perfluoroalkyl,
- 3) oxo,
- 4) OH,
- 5) halo,
- 6) CN,
- 7) (C<sub>2</sub>-C<sub>10</sub>)alkenyl,
- 8) (C<sub>2</sub>-C<sub>10</sub>)alkynyl,
- 9) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
- 10) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-aryl,
- 11) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-heterocyclyl,
- 12) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-N(R<sup>b</sup>)<sub>2</sub>,
- 13) C(O)R<sup>a</sup>,
- 14) (C<sub>0</sub>-C<sub>6</sub>)alkylene-CO<sub>2</sub>R<sup>a</sup>,
- 15) C(O)H,
- 16) (C<sub>0</sub>-C<sub>6</sub>)alkylene-CO<sub>2</sub>H,
- 17) C(O)N(R<sup>b</sup>)<sub>2</sub>,
- 18) S(O)<sub>m</sub>R<sup>a</sup>, and
- 19) S(O)<sub>2</sub>N(R<sup>b</sup>)<sub>2</sub>;

said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, and heterocyclyl is optionally substituted with up to three substituents selected from R<sup>b</sup>, OH, (C<sub>1</sub>-C<sub>6</sub>)alkoxy, halogen, CO<sub>2</sub>H, CN, O(C=O)C<sub>1</sub>-C<sub>6</sub> alkyl, oxo, and N(R<sup>b</sup>)<sub>2</sub>;

R<sup>12</sup> and R<sup>13</sup> are independently selected from:

- 1) H,
- 2) (C=O)O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 3) (C=O)O<sub>b</sub>C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
- 4) (C=O)Obaryl,
- 5) (C=O)Obheterocyclyl,
- 6) C<sub>1</sub>-C<sub>10</sub> alkyl,

- 7) aryl,
- 8) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 9) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 10) heterocyclyl,
- 11) C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
- 12) SO<sub>2</sub>R<sup>a</sup>, and
- 13) (C=O)NR<sup>b</sup><sub>2</sub>,

said alkyl, cycloalkyl, aryl, heterocyclyl, alkenyl, and alkynyl is optionally substituted with one, two or three substituents selected from R<sup>11</sup>, or

R<sup>12</sup> and R<sup>13</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>;

R<sup>a</sup> is independently selected from: (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, aryl, and heterocyclyl;

R<sup>b</sup> is independently selected from: H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, (C=O)OC<sub>1</sub>-C<sub>6</sub> alkyl, (C=O)C<sub>1</sub>-C<sub>6</sub> alkyl or S(O)<sub>2</sub>R<sup>a</sup>; and

R<sup>c</sup> and R<sup>c'</sup> are independently selected from: H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl and (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl or

R<sup>c</sup> and R<sup>c'</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>.

7. (Currently amended) The compound according to Claim 6 or a pharmaceutically acceptable salt or stereoisomer thereof, wherein:

R<sup>1</sup> is selected from:

- 1) (C=O)NR<sup>c</sup>RC',
- 2) ~~—SO<sub>2</sub>NR<sup>e</sup>Re'~~, and
- 3) ~~—SO<sub>2</sub>C<sub>1</sub>-C<sub>10</sub>-alkyl,~~

~~said alkyl, is optionally substituted with one, two or three substituents selected from R<sup>10</sup>;~~

R<sup>2</sup> is selected from:

- 1) aryl, and
- 2) ~~heteroaryl,~~

~~said aryl and heteroaryl is optionally substituted with one or more substituents selected from R<sup>10</sup>;~~

R<sup>3</sup>, R<sup>4</sup> and R<sup>8</sup> are independently selected from:

- 1) H, and
- 2) C<sub>1</sub>-C<sub>10</sub> alkyl,

~~said alkyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;~~

R<sup>6</sup> is selected from:

- 1) aryl, and
- 2) ~~heterocyclyl,~~

~~said alkyl, aryl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>, and.~~

~~R<sup>10</sup>, R<sup>11</sup>, R<sup>12</sup>, R<sup>13</sup>, Ra, Rb, Re and Re' are as described immediately above.~~

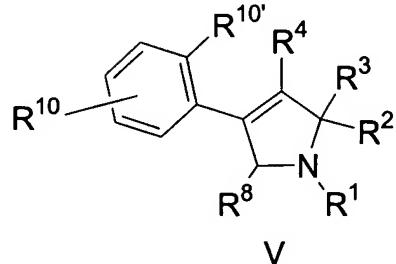
8. (Currently amended) The compound according to Claim 7 or a pharmaceutically acceptable salt or stereoisomer thereof, wherein:

R<sup>2</sup> and R<sup>6</sup> are independently selected from ~~are phenyl or pyridyl~~, optionally substituted with one or two substituents selected from R<sup>10</sup>, and.

~~R<sup>1</sup>, R<sup>3</sup>, R<sup>4</sup> and R<sup>8</sup> are as described in Claim 7.~~

9. (Cancelled)

10. (Currently amended) The compound according to Claim 5 of the Formula V,



wherein:

a is 0 or 1;  
b is 0 or 1;  
m is 0, 1, or 2;  
r is 0 or 1;  
s is 0 or 1;

R<sup>1</sup> is selected from:

- 1) (C=O)C<sub>1</sub>-C<sub>10</sub>-alkyl,
- 2) (C=O)aryl,
- 3) (C=O)C<sub>3</sub>-C<sub>8</sub>-cycloalkyl,
- 4) (C=O)heterocyclyl,
- 5) (C=O)NR<sup>c</sup>RC',
- 6) (C=S)NR<sup>e</sup>RE',
- 7) SO<sub>2</sub>NR<sup>e</sup>RE',
- 8) SO<sub>2</sub>C<sub>1</sub>-C<sub>10</sub>-alkyl,
- 9) SO<sub>2</sub>-aryl, and
- 10) SO<sub>2</sub>-heterocyclyl,

said alkyl, aryl, cycloalkyl, and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>2</sup> is selected from:

- 1) aryl,
- 2) ~~C<sub>1</sub>-C<sub>6</sub> aralkyl,~~
- 3) ~~C<sub>3</sub>-C<sub>8</sub> cycloalkyl, and~~
- 4) ~~heterocyclyl,~~

said aryl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>3</sup>, R<sup>4</sup> and R<sup>8</sup> are independently selected from:

- 1) H, and
- 2) C<sub>1</sub>-C<sub>10</sub> alkyl, and
- 3) ~~C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,~~

said alkyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>10</sup> is independently selected from:

- 1) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 2) (C=O)<sub>a</sub>O<sub>b</sub>aryl,
- 3) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 4) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 5) (C=O)<sub>a</sub>O<sub>b</sub> heterocyclyl,
- 6) CO<sub>2</sub>H,
- 7) halo,
- 8) CN,
- 9) OH,
- 10) O<sub>b</sub>C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,
- 11) O<sub>a</sub>(C=O)<sub>b</sub>NR<sup>12</sup>R<sup>13</sup>,
- 12) S(O)<sub>m</sub>R<sup>a</sup>,
- 13) S(O)<sub>2</sub>NR<sup>12</sup>R<sup>13</sup>,
- 14) ~~oxo,~~
- 15) CHO, and
- 16) ~~(N=O)R<sup>12</sup>R<sup>13</sup>, or~~

17)  $(C=O)_aObC_3-C_8$  cycloalkyl,  
said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one, two or three substituents selected from R<sup>11</sup>;  
R<sup>10'</sup> is halogen;

R<sup>11</sup> is selected from:

- 1)  $(C=O)_rOs(C_1-C_{10})$ alkyl,
- 2)  $O_f(C_1-C_3)$ perfluoroalkyl,
- 3) oxo,
- 4) OH,
- 5) halo,
- 6) CN,
- 7)  $(C_2-C_{10})$ alkenyl,
- 8)  $(C_2-C_{10})$ alkynyl,
- 9)  $(C=O)_rOs(C_3-C_6)$ cycloalkyl,
- 10)  $(C=O)_rOs(C_0-C_6)$ alkylene-aryl,
- 11)  $(C=O)_rOs(C_0-C_6)$ alkylene-heterocyclyl,
- 12)  $(C=O)_rOs(C_0-C_6)$ alkylene-N(R<sup>b</sup>)<sub>2</sub>,
- 13) C(O)R<sup>a</sup>,
- 14)  $(C_0-C_6)$ alkylene-CO<sub>2</sub>R<sup>a</sup>,
- 15) C(O)H,
- 16)  $(C_0-C_6)$ alkylene-CO<sub>2</sub>H,
- 17) C(O)N(R<sup>b</sup>)<sub>2</sub>,
- 18) S(O)<sub>m</sub>R<sup>a</sup>, and
- 19) S(O)<sub>2</sub>N(R<sup>b</sup>)<sub>2</sub>;

said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, alkylene and heterocyclyl is optionally substituted with up to three substituents selected from R<sup>b</sup>, OH, (C<sub>1</sub>-C<sub>6</sub>)alkoxy, halogen, CO<sub>2</sub>H, CN, O(C=O)C<sub>1</sub>-C<sub>6</sub> alkyl, oxo, and N(R<sup>b</sup>)<sub>2</sub>;

R<sup>12</sup> and R<sup>13</sup> are independently selected from:

- 1) H,
- 2)  $(C=O)ObC_1-C_{10}$  alkyl,

- 3)  $(C=O)ObC_3\text{-}C_8$  cycloalkyl,
- 4)  $(C=O)Ob$ aryl,
- 5)  $(C=O)Ob$ heterocyclyl,
- 6)  $C_1\text{-}C_{10}$  alkyl,
- 7) aryl,
- 8)  $C_2\text{-}C_{10}$  alkenyl,
- 9)  $C_2\text{-}C_{10}$  alkynyl,
- 10) heterocyclyl,
- 11)  $C_3\text{-}C_8$  cycloalkyl,
- 12)  $SO_2R^a$ , and
- 13)  $(C=O)NR^b_2$ ,

said alkyl, cycloalkyl, aryl, heterocyclyl, alkenyl, and alkynyl is optionally substituted with one, two or three substituents selected from R<sup>11</sup>, or

R<sup>12</sup> and R<sup>13</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>;

R<sup>a</sup> is independently selected from:  $(C_1\text{-}C_6)$ alkyl,  $(C_3\text{-}C_6)$ cycloalkyl, aryl, and heterocyclyl;

R<sup>b</sup> is independently selected from: H,  $(C_1\text{-}C_6)$ alkyl, aryl, heterocyclyl,  $(C_3\text{-}C_6)$ cycloalkyl,  $(C=O)OC_1\text{-}C_6$  alkyl,  $(C=O)C_1\text{-}C_6$  alkyl or  $S(O)_2R^a$ ; and

R<sup>c</sup> and R<sup>c'</sup> are independently selected from: H,  $(C_1\text{-}C_6)$ alkyl, aryl, heterocyclyl and  $(C_3\text{-}C_6)$ cycloalkyl or

R<sup>c</sup> and R<sup>c'</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said

monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R11.

11. (Currently amended) A compound selected from:

4-(2-chloro-5-fluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(+)-4-(2,5-difluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(-)-4-(2,5-difluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(5-chloro-2-fluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2-fluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2-fluoro-5-methylphenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(5-bromo-2-fluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4- {[4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}morpholine;

4- {[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}morpholine;

N,N-dimethyl-2,4-diphenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

~~3-[2-fluoro-5-(trifluoromethyl)phenyl]-N,N-dimethyl-5-phenyl-2,3-dihydro-1H-pyrrole-1-carboxamide;~~

2-(3-fluorophenyl)-4-(2,5-difluorophenyl)- N,N-dimethyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(4-fluorophenyl)-N,N-dimethyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(2-fluorophenyl)-N,N-dimethyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

2-(3-bromophenyl)-4-(2,5-difluorophenyl)- N,N-dimethyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

2-(3-aminophenyl)-4-(2,5-difluorophenyl)- N,N-dimethyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-methylphenyl)-N,N-dimethyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(5-chloro-2-fluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

~~4 (2,5 difluorophenyl) 1 (methylsulfonyl) 2 phenyl 2,5 dihydro 1H pyrrole;~~

~~4 (2,5 difluorophenyl) 1 (ethylsulfonyl) 2 phenyl 2,5 dihydro 1H pyrrole;~~

~~4 (2,5 difluorophenyl) 2 phenyl 1 (propylsulfonyl) 2,5 dihydro 1H pyrrole;~~

~~4 (2,5 difluorophenyl) 1 (isopropylsulfonyl) 2 phenyl 2,5 dihydro 1H pyrrole;~~

~~4 (5 chloro 2 fluorophenyl) 1 (methylsulfonyl) 2 phenyl 2,5 dihydro 1H pyrrole;~~

~~4 (5 chloro 2 fluorophenyl) 1 (isopropylsulfonyl) 2 phenyl 2,5 dihydro 1H pyrrole;~~

~~4 (2 fluoro 5 methylphenyl) 1 (isopropylsulfonyl) 2 phenyl 2,5 dihydro 1H pyrrole;~~

~~2 {[4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl]sulfonyl} ethanamine;~~

~~2 {[4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl]sulfonyl} N,N-dimethyl ethanamine;~~

~~1-acetyl-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~4-(2-chloro-5-fluorophenyl)-1-pivaloyl-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~4-(2-chloro-5-fluorophenyl)-1-isobutyryl-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~4-(2,5-difluorophenyl)-1-(2,2-dimethylpropanoyl)-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-methyl-1-oxopropan-2-ol;~~

~~1-[4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-methyl-1-oxopropan-2-ol;~~

~~1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-methyl-1-oxopropan-2-amine;~~

~~4-(2-fluoro-5-isocyanophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

~~4-(2,5-difluorophenyl)-2-phenyl-1-(trifluoroacetyl)-2,5-dihydro-1H-pyrrole;~~

~~4-(5-chloro-2-fluorophenyl)-2-phenyl-1-(trifluoroacetyl)-2,5-dihydro-1H-pyrrole;~~

~~(1S)-1-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl]-2-methylpropylamine;~~

~~(1R)-1-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl]-2-methylpropylamine;~~

~~4-(2,5-difluorophenyl)-2-phenyl-1-L-prolyl-2,5-dihydro-1H-pyrrole;~~

~~4-(2,5-difluorophenyl)-2-phenyl-1-D-prolyl-2,5-dihydro-1H-pyrrole;~~

(4R)-4-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl]-1,3-thiazolidine;

methyl (3S)-3-amino-4-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-4-oxobutanoate;

(4S)-4-amino-5-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-5-oxopentanamide;

(1S)-1-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl]-3-(methylthio)propylamine;

(1S)-1-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl]-3-(methylsulfonyl)propylamine;

(2S)-2-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl]piperidine;

(1S)-1-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl]pentylamine;

(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-(thien-2-ylmethyl)ethylamine;

4-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl]-1,1-dioxidotetrahydro-2H-thiopyran-4-ylamine;

(2S)-1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-methyl-1-oxopropan-2-amine;

(1S)-1-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl]propylamine;

(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-phenylethanamine;

(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-phenylethanamine;

(4S)-4-amino-5-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-5-exopentanamide

3-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-exopropan-1-amine;

(1S,2S)-1-([4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl)-2-methylbutylamine;

(1S)-1-([4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl)butylamine;

(1S)-1-cyclopropyl-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethanamine;

1-([4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl)cyclopropanamine;

1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-exopropan-2-amine;

(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-methyl-2-exoethylamine;

(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-(pyridin-2-ylmethyl)ethylamine;

(1S)-1-cyclohexyl-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethanamine;

(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-(4-iodobenzyl)-2-exoethylamine;

(1S)-1-benzyl-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethylamine;

4-((2S)-2-amino-3-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-exopropyl)phenol;

(*3S*) ~~3 {[4 (2,5-difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl]carbonyl} 1,2,3,4-tetrahydroisoquinoline;~~

(*1S*) ~~1 {[4 (2,5-difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl]carbonyl} 3 phenylpropylamine;~~

(*1S*) ~~1 {[4 (2,5-difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl]carbonyl} 3 methylbutylamine;~~

(*1S*) ~~2 [4 (2,5-difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl] 2 oxo 1 (pyridin 3 ylmethyl)ethylamine;~~

~~1 [(2*S*) azetidin 2 ylcarbonyl] 4 (2,5-difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrole;~~

(*3S*) ~~3 amino 4 [4 (2,5-difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl] 4 oxobutanamide; 4 (2,5-difluorophenyl) 1 [(2*S*) 2,5 dihydro 1H pyrrol 2 ylcarbonyl] 2 phenyl 2,5 dihydro 1H pyrrole;~~

~~4 (2,5-difluorophenyl) 1 [(2-methylazetidin 2 yl)carbonyl] 2 phenyl 2,5 dihydro 1H pyrrole;~~

(*1S*) ~~1 {[4 (2,5-difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl]carbonyl} 2,2 dimethylpropylamine;~~

~~methyl (4*S*) 4 amino 5 [4 (2,5-difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl] 5 exopentanoate;~~

~~4 (2,5-difluorophenyl) 2 phenyl 1 {[*(2S,3S)* 2 phenylpyrrolidin 3 yl]carbonyl} 2,5 dihydro 1H pyrrole;~~

~~4 (2,5-difluorophenyl) 2 phenyl 1 [(5 phenylpyrrolidin 3 yl)carbonyl] 2,5 dihydro 1H pyrrole;~~

(2S)-2-amino-3-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-oxopropan-1-ol;

(2R,3S)-3-amino-4-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-4-oxobutan-2-ol;

(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-(methoxymethyl)-2-exoethylamine;

4-(2,5-difluorophenyl)-2-phenyl-1-(pyrrolidin-3-ylcarbonyl)-2,5-dihydro-1H-pyrole;

4-(2,5-difluorophenyl)-2-phenyl-1-[(3-phenylpyrrolidin-3-yl)acetyl]-2,5-dihydro-1H-pyrole;

(1S)-1-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl]-3,3-difluoropropylamine;

(1S)-3-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-oxo-1-phenylpropan-1-amine;

4-(2,5-difluorophenyl)-2-phenyl-1-[(4S)-4-phenyl-L-prolyl]-2,5-dihydro-1H-pyrole;

1-{2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl}cyclohexanamine;

2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exethanamine;

4-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl]piperidin-4-amine;

(1S,3R)-3-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl]cyclopentanamine;

(1R,4S)-4-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl]cyclopent-2-en-1-amine;

(1S,4R) 4 {[4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl]carbonyl}cyclopent 2 en 1 amine;

(1S) 1 {[4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl]carbonyl}but 3 ynlyamine;

(1R) 3 [4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl] 3 oxo 1 phenylpropan 1 amine;

3 {[4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl]carbonyl} 2 phenylpiperidine;

(1S) 1 {[4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl]carbonyl}but 3 enylamine;

(2S) 3 [4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl] 2 (methylamine) 3 exopropan 1 ol;

(3R,5S) 5 {[4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl]carbonyl}pyrrolidin 3 ol;

(1S) 2 [4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl] 2 oxo 1 (1,3 thiazol 4 ylmethyl)ethylamine;

(1R) 1 {[4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl]carbonyl}but 3 enylamine;

(2S) 1 [4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl] N,3 dimethyl 1 exobutan 2 amine;

(2S) 1 [4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl] N,4 dimethyl 1 exopentan 2 amine;

(1S) 2 [4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl] 1 [(1 methyl 1H imidazol 4 yl)methyl] 2 oxoethylamine;

4 (2,5 difluorophenyl) 1 (N 6 formyl L lysyl) 2 phenyl 2,5 dihydro 1H pyrrole;

(*2S,3S*)-1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-N,3-dimethyl-1-exo-pentan-2-amine;

(*1S*)-1-(cyclohexylmethyl)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-exo-ethylamine;

(*1S*)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-1-(1*H*-indol-3-ylmethyl)-2-exo-ethylamine;

(*1S*)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-1-(isocyanoethyl)-2-exo-ethylamine;

(*1S*)-1-([4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]carbonyl)-3,3-dimethylbutylamine;

1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2,3-dimethyl-1-exobutan-2-amine;

1-([4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]carbonyl)cyclohexanamine;

1-([4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]carbonyl)cyclopentanamine;

(*1S*)-3-(benzylxy)-1-([4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]carbonyl)propylamine;

1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2,3-dimethyl-1-exobutan-2-amine;

1-([4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]carbonyl)cyclopent-3-en-1-amine;

(*S*) 1-cyclopentyl 2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-exoethanamine;

4-(2,5-difluorophenyl)-1-(2-methylpropyl)-2-phenyl-2,5-dihydro-1*H*-pyrrole;

1-[4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-methyl-1-exopropan-2-amine;

(*S*) 1-[[4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]carbonyl]-2-methylpropylamine;

(*S*) 2-[4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-1-cyclopropyl-2-exoethanamine;

(*S,2S*) 1-[[4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]carbonyl]-2-methylbutylamine;

(*S*) 1-[[4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]carbonyl]pentylamine;

(*S*) 1-[[4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]carbonyl]-3,3-dimethylbutylamine;

(*S*) 1-[[*(2S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]carbonyl]-2,2-dimethylpropylamine;

(*S*) 1-[[*(2S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]carbonyl]-2-methylpropylamine;

(*S*) 1-cyclohexyl-2-[(*2S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-exoethanamine;

(<sup>1S</sup>) 1 [(<sup>2S</sup>) 4 (2,5-difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl] carbonyl but 3-enylamine;

(<sup>1S</sup>) 1 [(<sup>2S</sup>) 4 (2,5-difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl] carbonyl but 3-ynylamine;

(<sup>1S</sup>) 1 benzyl 2 [<sup>(2S)</sup> 4 (2,5-difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl] 2-exoethylamine;

(<sup>1S</sup>) 1 cyclopropyl 2 [<sup>(2S)</sup> 4 (2,5-difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl] 2-exoethanamine;

1 [(<sup>2S</sup>) 4 (2,5-difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl] 2-methyl 1 exopropan 2-amine;

(<sup>1S</sup>) 1 [(<sup>2S</sup>) 4 (5 chloro 2 fluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl] carbonyl 2,2-dimethylpropylamine;

(<sup>1S</sup>) 1 [(<sup>2S</sup>) 4 (2,5-difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl] carbonyl pentylamine;

(<sup>1S</sup>) 1 [(<sup>2S</sup>) 4 (2,5-difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl] carbonyl 3-methylbutylamine;

(<sup>1S</sup>) 1 [(<sup>2S</sup>) 4 (2,5-difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl] carbonyl 3,3-dimethylbutylamine;

1 cyclopropyl 3 [<sup>(2S)</sup> 4 (2,5-difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl] 3-exopropan 1 amine;

(<sup>1S</sup>) 2 [<sup>(2S)</sup> 4 (5 chloro 2 fluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol 1 yl] 1 cyclopropyl 2-exoethanamine;

(*1S*) 1  $\{[(2S)-4-(5\text{-chloro-2-fluorophenyl})-2\text{-phenyl}-2,5\text{-dihydro-1H-pyrrol-1-yl]carbonyl}\} - 2\text{-methylpropylamine;}$

(*1S,2S*) 1  $\{[(2S)-4-(2,5\text{-difluorophenyl})-2\text{-phenyl}-2,5\text{-dihydro-1H-pyrrol-1-yl]carbonyl}\} - 2\text{-methylbutylamine;}$

4-(5-chloro-2-fluorophenyl)-2-(3-hydroxyphenyl)-1-(2-methylalanyl)-2,5-dihydro-1H-pyrrole;

(*2S*) 4-(5-chloro-2-fluorophenyl)-2-(3-hydroxyphenyl)-1-(2-methylalanyl)-2,5-dihydro-1H-pyrrole;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-1-L-valyl-2,5-dihydro-1H-pyrrole;

4-(5-chloro-2-fluorophenyl)-2-(3-hydroxyphenyl)-1-L-valyl-2,5-dihydro-1H-pyrrole;

(*2S*) 4-(5-chloro-2-fluorophenyl)-2-(3-hydroxyphenyl)-1-L-valyl-2,5-dihydro-1H-pyrrole;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-1-(2-methylalanyl)-2,5-dihydro-1H-pyrrole;

3-[1-[(*2S*)-2-amino-2-cyclopropylethanoyl]-4-(5-chloro-2-fluorophenyl)-2,5-dihydro-1H-pyrrol-2-yl]phenol;

4-(5-chloro-2-fluorophenyl)-2-(3-hydroxyphenyl)-1-L-isoleucyl-2,5-dihydro-1H-pyrrole;

4-(5-chloro-2-fluorophenyl)-2-(3-hydroxyphenyl)-1-L-norleucyl-2,5-dihydro-1H-pyrrole;

(*2S*) 4-(5-chloro-2-fluorophenyl)-2-(3-hydroxyphenyl)-1-(3-methyl-L-valyl)-2,5-dihydro-1H-pyrrole;

(*2S*)-4-(2,5-Difluorophenyl)-N-methyl-2-phenyl-N-piperidin-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(2,5-Difluorophenyl)-N-methyl-2-phenyl-N-(piperidin-4-ylmethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(5-Chloro-2-fluorophenyl)-N-methyl-2-phenyl-N-[(3R)-pyrrolidin-3-yl]-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(5-chloro-2-fluorophenyl)-N-methyl-2-phenyl-N-piperidin-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(5-chloro-2-fluorophenyl)-N-methyl-2-phenyl-N-[(3S)-pyrrolidin-3-yl]-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-pyrrolidin-3-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(2,5-difluorophenyl)-N-methyl-N-[(3S)-1-methylpyrrolidin-3-yl]-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(2,5-difluorophenyl)-N-methyl-N-[(3R)-1-methylpyrrolidin-3-yl]-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(2,5-difluorophenyl)-N-methyl-N-[(1-methyl-5-oxopyrrolidin-2-yl)methyl]-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(2,5-difluorophenyl)-N-(1,3-dioxolan-2-ylmethyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-tetrahydrofuran-3-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-N-(1-allylpiperidin-4-yl)-4-(2,5-difluorophenyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

allyl 4-[{[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}(methyl)amino]piperidine-1-carboxylate;

allyl 4-{[{(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}(methyl)amino]methyl}piperidine-1-carboxylate;

(2S)-4-(2,5-difluorophenyl)-N-methyl-N-(1-methylpiperidin-4-yl)-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-Difluorophenyl)-N-methyl-N-[(1-methylpiperidin-3-yl)methyl]-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-(pyridin-3-ylmethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;

N-benzyl-4-(2,5-difluorophenyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-methyl-N-[(1-methyl-1H-pyrazol-4-yl)methyl]-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-[2-(dimethylamino)ethyl]-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-(2-hydroxyethyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-isobutyl-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-(2-pyridin-2-ylethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-(2-methoxyethyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-(2,3-dihydroxypropyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-(2-phenylethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-phenyl-N-propyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(5-chloro-2-fluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-piperidin-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-piperidin-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;

~~1-Acetyl 4-(2,5-difluorophenyl)-2-methyl-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~(2S)-1-[4-(2,5-difluorophenyl)-2-methyl-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]3-methyl-1-exobutan-2-amine;~~

(2S)-4-(2,5-difluorophenyl)-N,N,2-trimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(5-chloro-2-fluorophenyl)-N,N,2-trimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(5-chloro-2-fluorophenyl)-2-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(5-chloro-2-fluorophenyl)-N-ethyl-2-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3,3-dimethyl-1-exobutan-2-ol;

(2S)-1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-methyl-1-exobutan-2-ol;

(2S,3S)-1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-methyl-1-exopentan-2-ol;

1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-exohexan-2-ol;

(2S)-1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-oxo-3-phenylpropan-2-ol;

(2S)-1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-4-methyl-1-exopentan-2-ol;

(1S)-1-cyclohexyl-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethanol;

(2S)-1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3,3-dimethyl-1-exobutan-2-ol;

N-1-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl)-N-2,N-2-dimethylglycinamide;

N-1-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl)-N-2-methylglycinamide;

N-1-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl)glycinamide;

~~N-1 {((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl}-2-methylalaninamide;~~

~~N-1 {((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl}-glycinamide;~~

~~N-1 {((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl}-N-2,N-2-dimethylglycinamide;~~

~~N-1 {((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl}-N-2,N-2-dimethylglycinamide, N-oxide;~~

~~N-1 {((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl}-2-methylalaninamide;~~

~~N-1 {((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl}-N-2,N-2-dimethylglycinamide n-oxide;~~

~~N-1 {((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl}-2-pyrrolidin-1-ylacetamide;~~

~~2-azetidin-1-yl N-1 {((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl}-2-acetamide;~~

~~N-1 {((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl}-2-morpholin-4-ylacetamide;~~

~~N-1 {((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl}-2-piperazin-1-ylacetamide;~~

~~N-1 {((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl}-2-(4-methylpiperazin-1-yl)acetamide;~~

~~2-azetidin-1-yl-N-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl)acetamide;~~

~~N-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl)-2-pyrrolidin-1-ylacetamide;~~

~~N-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl)-2-piperidin-1-ylacetamide;~~

~~N-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl)-2-morpholin-4-ylacetamide;~~

~~N-1-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl)-N-2-(2-hydroxyethyl)glycinamide;~~

~~N-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl)-2-(4-methylpiperazin-1-yl)acetamide;~~

~~N-1-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl)-N-2-isopropylglycinamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl)acetamide;~~

~~N-1-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl)-N-2-ethylglycinamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl)-2-hydroxyacetamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl)piperazine-1-carboxamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-N'-piperidin-4-ylurea;~~

~~4-amino-N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)piperidine-1-carboxamide;~~

~~N-(2-aminoethyl)-N'-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)urea;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-N'-(3-morpholin-4-ylpropyl)urea;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-N'[2-(dimethylamino)ethyl]urea;~~

~~2-azetidin-1-yl-N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)ethanesulfonamide~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-(isopropylamino)ethanesulfonamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-pyrrolidin-1-ylethanesulfonamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-morpholin-4-ylethanesulfonamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-piperazin-1-ylethanesulfonamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-(4-methylpiperazin-1-yl)ethanesulfonamide;~~

N-(tert-butyl)-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]acetamide;

2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-isopropylacetamide;

(2S)-1-(2-azetidin-1-yl-2-oxoethyl)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrole;

(2S)-4-(2,5-difluorophenyl)-1-(2-oxo-2-pyrrolidin-1-ylethyl)-2-phenyl-2,5-dihydro-1H-pyrrole;

4-{[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]acetyl}morpholine;

1-{[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]acetyl}piperazine;

1-{[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]acetyl}-4-methylpiperazine;

2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-isopropylbutanamide;

4-{2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]butanoyl}morpholine;

2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-ethylacetamide;

N-cyclobutyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]acetamide;

2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-ethylpropanamide;

N-cyclobutyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanamide;

2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-methylpropanamide;

2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-isopropylpropanamide;

N-(tert-butyl)-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanamide;

4-{(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanoyl}morpholine;

(3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-ethyl-2,2-dimethyl-4-oxobutanamide;

(3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2,2-dimethyl-4-oxo-N-piperidin-4-ylbutanamide;

(3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2,2-dimethyl-4-oxobutanoic acid;

(3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,N,2,2-tetramethyl-4-oxobutanamide;

(1S)-1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2,2-dimethyl-3-oxo-3-piperazin-1-ylpropylamine;

(3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-isopropyl-2,2-dimethyl-4-oxobutanamide;

(3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,2,2-trimethyl-4-oxobutanamide;

(3R)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,N,2,2-tetramethyl-4-oxobutanamide;

(3R)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2,2-dimethyl-4-oxobutanoic acid;

(1R)-1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2,2-dimethyl-3-oxo-3-piperazin-1-ylpropylamine;

~~2-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl)amino-N-ethylacetamide;~~

~~2-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl)amino-N-methylacetamide;~~

~~2-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl)amino-N,N-dimethylacetamide;~~

~~2-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl)amino-N-methyl-N-ethylacetamide;~~

~~2-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl)amino-N-methylacetamide;~~

~~2-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl)amino-N-ethylacetamide;~~

~~2-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl)amino-N,N-dimethylacetamide;~~

~~2-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl)amino-N-isopropylacetamide;~~

~~2-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl)amino-N-ethyl-N-methylacetamide;~~

~~2-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl)amino-N,N-diethylacetamide;~~

~~(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exo-N-(2-exo-2-pyrrolidin-1-ylethyl)ethanamine;~~

(*S*) 1 cyclopropyl 2 [(*S*) 4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1*H* pyrrol 1 yl] N (2 morpholin 4 yl 2 oxoethyl) 2 oxoethanamine;

1 [{((*S*) 1 cyclopropyl 2 [(*S*) 4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1*H* pyrrol 1 yl] 2 oxoethyl} amino]acetyl]piperidin 4 ol;

(*S*) 1 cyclopropyl 2 [(*S*) 4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1*H* pyrrol 1 yl] N [2 (4 methylpiperazin 1 yl) 2 oxoethyl] 2 oxoethanamine;

(*S*) N (2 azetidin 1 yl 2 oxoethyl) 1 cyclopropyl 2 [(*S*) 4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1*H* pyrrol 1 yl] 2 oxoethanamine;

(*S*) 1 cyclopropyl 2 [(*S*) 4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1*H* pyrrol 1 yl] N [2 (1,1 dioxidothiomorpholin 4 yl) 2 oxoethyl] 2 oxoethanamine;

(*S*) N [2 (4 acetyl piperazin 1 yl) 2 oxoethyl] 1 cyclopropyl 2 [(*S*) 4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1*H* pyrrol 1 yl] 2 oxoethanamine;

(*S*) 1 tert butyl 2 [(*S*) 4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1*H* pyrrol 1 yl] N (2 morpholin 4 yl 2 oxoethyl) 2 oxoethanamine;

(*S*) 1 tert butyl 2 [(*S*) 4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1*H* pyrrol 1 yl] 2 o xo N (2 o xo 2 pyrrolidin 1 ylethyl)ethanamine;

2-({(*S*)-1-tert-butyl-2-[*(S)*-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-oxoethyl} amino)-N-isopropylacetamide;

2-(dimethylamino)ethyl (*S*)-1-cyclopropyl-2-[*(S)*-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-oxoethylcarbamate;

1-methylpiperidin-4-yl (*S*)-1-cyclopropyl-2-[*(S)*-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-oxoethylcarbamate;

(2S)-4-cyclopropyl-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-cyclopentyl-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl-4-methylpiperazine-1-carboxylate;

1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl-2-morpholin-4-ylethylcarbamate;

N-[(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl]oxy)carbonyl]glycine;

(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl-1-methylpiperidin-4-ylcarbamate;

(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethylmethyl(1-methylpiperidin-4-yl)carbamate;

(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl-4-dimethylamino)piperidine-1-carboxylate;

tert-butyl (2S)-4-(2-chloro-5-fluoropyrimidin-4-yl)-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxylate;

(2S)-4-(5-fluoro-2-methylpyrimidin-4-yl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(2-chloro-5-fluoropyrimidin-4-yl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(4-chloro-5-methylpyrimidin-2-yl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(6-chloropyrimidin-4-yl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(2-chloropyrimidin-4-yl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-N,N-dimethyl-4-(4-methylpyridin-3-yl)-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-N,N-dimethyl-2-phenyl-4-(1,3-thiazol-2-yl)-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-N,N-dimethyl-2-phenyl-4-(1,3-thiazol-4-yl)-2,5-dihydro-1H-pyrrole-1-carboxamide;  
4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N,N-dimethyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N,N-dimethyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(2,5-difluorophenyl)-N-(2-hydroxyethyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

N-{[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-N-methyl-beta-alanine;

methyl N-{[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-N-methyl-beta-alaninate;

4-{[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]acetyl}morpholin-4-ium;

~~3-[2S)-4-(2,5-difluorophenyl)-1-(2-hydroxy-2-methylpropanoyl)-2,5-dihydro-1H-pyrrol-2-yl]phenol;~~

~~4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N,N-dimethyl-2,5-dihydro-1H-pyrrole-1-sulfonamide;~~

~~3-[4-(2,5-difluorophenyl)-1-(methylsulfonyl)-2,5-dihydro-1H-pyrrol-2-yl]phenol;~~

~~3-[4-(2,5-difluorophenyl)-1-(morpholin-4-ylcarbonyl)-2,5-dihydro-1H-pyrrol-2-yl]phenol;~~

~~3-[4-(2,5-difluorophenyl)-1-(2,2-dimethylpropanoyl)-2,5-dihydro-1H-pyrrol-2-yl]phenol;~~

~~(2S)-4-(2,5-difluorophenyl)-1-[(methylsulfonyl)acetyl]-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~(2S)-4-(2,5-difluorophenyl)-2-phenyl-1-[(phenylsulfonyl)acetyl]-2,5-dihydro-1H-pyrrole;~~

~~3-[{(2S)-1-[(2S)-2-cyclopropyl-2-hydroxyethanoyl]}-4-(2,5-difluorophenyl)-2,5-dihydro-1H-pyrrol-2-yl]phenol;~~

~~3-((2S)-4-(2,5-difluorophenyl)-1-[(2S)-2-hydroxy-3,3-dimethylbutanoyl]-2,5-dihydro-1H-pyrrol-2-yl]phenol;~~

~~(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethanol;~~

~~(2S)-4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-tetrahydrofuran-3-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

~~(2S)-4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-(2-methoxyethyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

~~(2S)-4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-tetrahydro-2H-pyran-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

~~3-[(2S)-4-(2,5-difluorophenyl)-1-(piperidin-1-ylcarbonyl)-2,5-dihydro-1H-pyrrol-2-yl]phenol;~~

4-(2,5-difluorophenyl)-N-[1-(2-fluoroethyl)piperidin-4-yl]-2-(3-hydroxyphenyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-[{[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}(methyl)amino]piperidinium trifluoroacetate;

2-[{[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}(methyl)amino]ethyl 4-methylpiperazine-1-carboxylate;

3-{4-(2,5-difluorophenyl)-1-[(4-methylpiperazin-1-yl)carbonyl]-2,5-dihydro-1H-pyrrol-2-yl}phenol;

2-[{[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}(methyl)amino]ethyl morpholine-4-carboxylate;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(1-methyl-1H-pyrazol-4-yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-(isoxazol-5-ylmethyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

2-[{[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}(methyl)amino]ethyl dimethylaminocarboxylate;

2-[{[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}(methyl)amino]ethyl piperidine-1-carboxylate;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[2-(2-oxopyrrolidin-1-yl)ethyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(5-oxo-4,5-dihydro-1H-1,2,4-triazol-3-yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-(tetrahydro-2H-pyran-4-ylmethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-{{[5-(methoxymethyl)-1H-pyrazol-3-yl]methyl}-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-(1,3-thiazol-4-ylmethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(4-methyl-1,2,5-oxadiazol-3-yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-(1,3-thiazol-2-ylmethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-(isoxazol-3-ylmethyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[2-(1H-1,2,4-triazol-1-yl)ethyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[2-(1H-pyrazol-1-yl)ethyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(1-methyl-5-oxopyrrolidin-2-yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-(1-isoxazol-3-ylethyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-(1,3-dioxolan-2-ylmethyl)-2-(3-hydroxyphenyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-(1,4-dioxan-2-ylmethyl)-2-(3-hydroxyphenyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(5-methyl-1,3,4-oxadiazol-2-yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[2-(methylsulfonyl)ethyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;

2-[{[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}(methyl)amino]ethanesulfonic acid;

~~2-hydroxyethyl (1S)-1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl-2,2-dimethylpropylcarbamate;~~

~~3-hydroxypropyl (1S)-1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl-2,2-dimethylpropylcarbamate;~~

~~2-hydroxyethyl ((1S)-1-isopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)carbamate;~~

~~2-hydroxyethyl ((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)carbamate;~~

~~4-hydroxybutyl (1S)-1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl-2,2-dimethylpropylcarbamate;~~

~~(2S)-4-(2,5-difluorophenyl)-1-[2-(methylsulfonyl)ethyl]-2-phenyl-2,5-dihydro-1H-pyrrole;~~

(2S)-4-(2,5-difluorophenyl)-1-[2-(ethylsulfonyl)ethyl]-2-phenyl-2,5-dihydro-1*H*-pyrrole;  
1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]pentan-3-one;  
4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]butan-2-one;  
4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-3-methylbutan-2-one;  
2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-*N,N*-dimethylethanedisulfonamide;  
3-[(2S)-4-(2,5-difluorophenyl)-1-[2-(methylsulfonyl)ethyl]-2,5-dihydro-1*H*-pyrrol-2-yl]phenol;  
methyl 3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]propanoate;  
(2S)-4-(2,5-difluorophenyl)-1-[2-(ethylsulfonyl)propyl]-2-phenyl-2,5-dihydro-1*H*-pyrrole;  
3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-*N*-methylpropanamide;  
3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-*N,N*-dimethylpropanamide;  
3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-*N,N*,2-trimethylpropanamide;  
4-{3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]propanoyl}morpholine;  
1-{3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]propanoyl}-4-(methylsulfonyl)piperazine;  
1-{3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]propanoyl}piperidin-4-ol;  
methyl 3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]propanoate;  
2-((1*S*)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-*exo*ethyl)oxy)-*N*-ethylacetamide;  
4-((1*S*)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-*exo*ethoxy)acetyl)morpholine;

~~2-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethoxy) N-(2-hydroxyethyl)acetamide;~~

~~1-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethoxy)acetyl) 4-methylpiperazine;~~

~~1-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethoxy)acetyl)piperazine;~~

~~2-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethoxy) N-piperidin-4-ylacetamide;~~

~~1-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethoxy)acetyl)piperidin-4-amine;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-3-morpholin-4-yl-3-oxopropyl-1-amine;~~

~~N<sup>3</sup>-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-N<sup>1</sup>,N<sup>1</sup>-dimethyl-β-alaninamide; and~~

~~((1S)-1-((2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl)-2,2-dimethylpropyl)(3-morpholin-4-yl-3-oxopropyl)amine;~~

or a pharmaceutically acceptable salt or stereoisomer thereof.

12. (Currently amended) The compound according to Claim 11 which is selected from:

~~(-)-4-(2,5-difluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

~~4-(5-chloro-2-fluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

~~4-(2,5-difluorophenyl)-1-(isopropylsulfonyl)-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-methyl-1-oxopropan-2-ol;~~

~~1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-methyl-1-oxopropan-2-amine;~~

~~(1S)-1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-2,2-dimethylpropylamine; and~~

~~(2S)-4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-piperidin-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

or a pharmaceutically acceptable salt or stereoisomer thereof.

13. (Currently amended) The compound according to Claim 11 which is selected from:

~~(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethanamine;~~

~~(2S)-4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-piperidin-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

~~(1S)-1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-2,2-dimethylpropylamine;~~

~~2-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl)amino)-N-ethylacetamide;~~

~~(2S)-4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N,N-dimethyl-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

~~(2S)-4-(2,5-difluorophenyl)-N-methyl-N-(1-methylpiperidin-4-yl)-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide; and~~

~~(2S)-4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-1-L-valyl-2,5-dihydro-1H-pyrrole~~

or a pharmaceutically acceptable salt or stereoisomer thereof.

14. (Currently amended) A compound which is:

~~(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethanamine~~

~~(2S)-4-(2,5-difluorophenyl)-1-[(methylsulfonyl)acetyl]-2-phenyl-2,5-dihydro-1H-pyrrole~~

~~(2S)-4-(2,5-difluorophenyl)-2-phenyl-1-[(phenylsulfonyl)acetyl]-2,5-dihydro-1H-pyrrole~~

~~3-[(2S)-1-[(2S)-2-cyclopropyl-2-hydroxyethanoyl]-4-(2,5-difluorophenyl)-2,5-dihydro-1H-pyrrol-2-yl]phenol~~

~~3-[(2S)-4-(2,5-difluorophenyl)-1-[(2S)-2-hydroxy-3,3-dimethylbutanoyl]-2,5-dihydro-1H-pyrrol-2-yl]phenol~~

~~(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethanol~~

~~(2S)-4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-tetrahydrofuran-3-yl-2,5-dihydro-1H-pyrrole-1-carboxamide~~

~~(2S)-4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-(2-methoxyethyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide~~

~~(2S)-4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-tetrahydro-2H-pyran-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide~~

~~3-[(2S)-4-(2,5-difluorophenyl)-1-(piperidin-1-ylcarbonyl)-2,5-dihydro-1H-pyrrol-2-yl]phenol~~

~~4-(2,5-difluorophenyl)-N-[1-(2-fluoroethyl)piperidin-4-yl]-2-(3-hydroxyphenyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide~~

4-[{[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}(methyl)amino]piperidinium trifluoroacetate

2-[{[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}(methyl)amino]ethyl 4-methylpiperazine-1-carboxylate

3-{4-(2,5-difluorophenyl)-1-[(4-methylpiperazin-1-yl)carbonyl]-2,5-dihydro-1H-pyrrol-2-yl}phenol

2-[{[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}(methyl)amino]ethyl morpholine-4-carboxylate

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(1-methyl-1H-pyrazol-4-yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-(isoxazol-5-ylmethyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide

2-[{[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}(methyl)amino]ethyl dimethylaminocarboxylate

2-[{[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}(methyl)amino]ethyl piperidine-1-carboxylate

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[2-(2-oxopyrrolidin-1-yl)ethyl]-2,5-dihydro-1H-pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(5-oxo-4,5-dihydro-1H-1,2,4-triazol-3-yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-(tetrahydro-2H-pyran-4-ylmethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N- {[5-(methoxymethyl)-1H-pyrazol-3-yl]methyl}-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-(1,3-thiazol-4-ylmethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(4-methyl-1,2,5-oxadiazol-3-yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-(1,3-thiazol-2-ylmethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-(isoxazol-3-ylmethyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[2-(1H-1,2,4-triazol-1-yl)ethyl]-2,5-dihydro-1H-pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[2-(1H-pyrazol-1-yl)ethyl]-2,5-dihydro-1H-pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(1-methyl-5-oxopyrrolidin-2-yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-(1-isoxazol-3-ylethyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-N-(1,3-dioxolan-2-ylmethyl)-2-(3-hydroxyphenyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-N-(1,4-dioxan-2-ylmethyl)-2-(3-hydroxyphenyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(5-methyl-1,3,4-oxadiazol-2-yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[2-(methylsulfonyl)ethyl]-2,5-dihydro-1H-pyrrole-1-carboxamide

2-[{[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}(methyl)amino]ethanesulfonic acid

~~2 hydroxyethyl (1S)-1 [(2S)-4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol-1-yl]carbonyl} 2,2 dimethylpropylecarbamate~~

~~3 hydroxypropyl (1S)-1 [(2S)-4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol-1-yl]carbonyl} 2,2 dimethylpropylecarbamate~~

~~2 hydroxyethyl ((1S)-1 isopropyl 2 [(2S)-4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol-1-yl] 2 oxoethyl} carbamate~~

~~2 hydroxyethyl ((1S)-1 cyclopropyl 2 [(2S)-4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol-1-yl] 2 oxoethyl} carbamate~~

~~4 hydroxybutyl (1S)-1 [(2S)-4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol-1-yl]carbonyl} 2,2 dimethylpropylecarbamate~~

~~(2S)-4 (2,5 difluorophenyl)-1 [2 (methylsulfonyl)ethyl] 2 phenyl 2,5 dihydro 1H pyrrole~~

~~(2S)-4 (2,5 difluorophenyl)-1 [2 (ethylsulfonyl)ethyl] 2 phenyl 2,5 dihydro 1H pyrrole~~

~~1 [(2S)-4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol-1-yl]pentan-3-one~~

~~4 [(2S)-4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol-1-yl]butan-2-one~~

~~4 [(2S)-4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol-1-yl] 3 methylbutan-2-one~~

~~2 [(2S)-4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol-1-yl] N,N-dimethylethanesulfonamide~~

~~3 [(2S)-4 (2,5 difluorophenyl) 1 [2 (methylsulfonyl)ethyl] 2,5 dihydro 1H pyrrol-2-yl}phenol~~

~~methyl 3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanoate~~

~~(2S)-4-(2,5-difluorophenyl)-1-[2-(ethylsulfonyl)propyl]-2-phenyl-2,5-dihydro-1H-pyrrole~~

~~3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-methylpropanamide~~

~~3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,N-dimethylpropanamide~~

~~3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,N,2-trimethylpropanamide~~

~~4-{3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanoyl}morpholine~~

~~1-{3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanoyl}-4-(methylsulfonyl)piperazine~~

~~1-{3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanoyl}piperidin-4-ol~~

~~methyl 3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanoate~~

~~2-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-  
exoethoxy)N-ethylacetamide~~

~~4-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-  
exoethoxy)acetyl)morpholine~~

~~2-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-  
exoethoxy)N-(2-hydroxyethyl)acetamide~~

~~1-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-  
exoethoxy)acetyl)-4-methylpiperazine~~

~~1-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-  
exoethoxy)acetyl)piperazine~~

~~2-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-  
exoethoxy)N-piperidin-4-ylacetamide~~

~~1-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-  
exoethoxy)acetyl)piperidin-4-amine~~

~~N<sup>[(1S)-1 cyclopropyl 2 [(2S)-4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol-1-yl] 2 exoethyl] 3 morpholin-4-yl 3 oxopropan-1-amine}</sup>~~

~~N<sup>3</sup>-[(1S)-1 cyclopropyl 2 [(2S)-4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol-1-yl] 2 exoethyl] N<sup>4</sup>,N<sup>4</sup>-dimethyl β alaninamide~~

~~((1S)-1 [(2S)-4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol-1-yl]carbonyl) 2,2-dimethylpropyl)(3 morpholin-4-yl 3 oxopropan-1-amine~~

or a pharmaceutically acceptable salt thereof.

15. (Currently amended) The compound according to Claim 12 which is the TFA salt of a compound selected from:

~~2-[(4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol-1-yl)sulfonyl] N,N-dimethylethanamine;~~

~~1-[4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol-1-yl]-2-methyl-1-oxopropan-2-amine;~~

~~4-(5 chloro-2-fluorophenyl) 2 phenyl 1-(trifluoroacetyl) 2,5 dihydro 1H pyrrole;~~

~~(1S)-1-[(4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol-1-yl]carbonyl) 2-methylpropylamine;~~

~~(1R)-1-[(4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol-1-yl]carbonyl) 2-methylpropylamine;~~

~~4-(2,5 difluorophenyl) 2 phenyl 1-L-prolyl 2,5 dihydro 1H pyrrole;~~

~~4-(2,5 difluorophenyl) 2 phenyl 1-D-prolyl 2,5 dihydro 1H pyrrole;~~

~~(4R)-4-[(4 (2,5 difluorophenyl) 2 phenyl 2,5 dihydro 1H pyrrol-1-yl]carbonyl) 1,3-thiazolidine;~~

~~methyl (3S)-3-amino-4-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-4-exobutanoate;~~

~~(4S)-4-amino-5-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-5-exopentanamide;~~

~~(1S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-3-(methylthio)propylamine;~~

~~(1S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-3-(methylsulfonyl)propylamine;~~

~~(2S)-2-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]piperidine;~~

~~(1S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]pentylamine;~~

~~(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-(thien-2-ylmethyl)ethylamine;~~

~~4-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-1,1-dioxidotetrahydro-2H-thiopyran-4-ylamine;~~

~~(2S)-1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-methyl-1-exopropan-2-amine;~~

~~(1S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]propylamine;~~

~~(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-phenylethanamine;~~

~~(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-phenylethanamine;~~

(4S)-4-amino-5-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-5-exo-pentanamide;

3-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-oxopropan-1-amine;

(1S,2S)-1-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl]-2-methylbutylamine;

(1S)-1-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl]butylamine;

(1S)-1-cyclopropyl-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethanamine;

1-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl]cyclopropanamine;

1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-oxopropan-2-amine;

(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-methyl-2-exoethylamine;

(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-(pyridin-2-ylmethyl)ethylamine;

(1S)-1-cyclohexyl-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethanamine;

(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-(4-iodobenzyl)-2-exoethylamine;

(1S)-1-benzyl-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethylamine;

4-((2S)-2-amino-3-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-exopropyl)phenol;

(3S)-3-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl]-1,2,3,4-tetrahydroisoquinoline;

(1S)-1-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl]-3-phenylpropylamine;

(1S)-1-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl]-3-methylbutylamine;

(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-(pyridin-3-ylmethyl)ethylamine;

1-[(2S)-azetidin-2-ylcarbonyl]-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrole;

(3S)-3-amino-4-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)-4-oxobutanamide];

4-(2,5-difluorophenyl)-1-[(2-methylazetidin-2-yl)carbonyl]-2-phenyl-2,5-dihydro-1H-pyrole;

(1S)-1-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl]-2,2-dimethylpropylamine;

methyl (4S)-4-amino-5-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)-5-oxopentanoate];

4-(2,5-difluorophenyl)-2-phenyl-1-[(2S,3S)-2-phenylpyrrolidin-3-yl]carbonyl]-2,5-dihydro-1H-pyrole;

4-(2,5-difluorophenyl)-2-phenyl-1-[(5-phenylpyrrolidin-3-yl)carbonyl]-2,5-dihydro-1H-pyrole;

(2S)-2-amino-3-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)-3-oxopropan-1-ol];

(2R,3S)-3-amino-4-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)-4-oxobut-2-enyl];

(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-(methoxymethyl)-2-  
exoeethylamine;

4-(2,5-difluorophenyl)-2-phenyl-1-(pyrrolidin-3-ylcarbonyl)-2,5-dihydro-1H-pyrrole;

4-(2,5-difluorophenyl)-2-phenyl-1-[(3-phenylpyrrolidin-3-yl)acetyl]-2,5-dihydro-1H-pyrrole;

(1S)-1-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl]-3,3-  
difluoropropylamine;

(1S)-3-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-oxo-1-phenylpropan-1-  
amine;

4-(2,5-difluorophenyl)-2-phenyl-1-[(4S)-4-phenyl-L-prolyl]-2,5-dihydro-1H-pyrrole;

1-{2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-  
exoeethyl}cyclohexanamine;

2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethanamine;

4-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl]piperidin-4-amine;

(1S,3R)-3-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-  
yl)carbonyl]cyclopentanamine;

(1R,4S)-4-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl]cyclopent-2-  
en-1-amine;

(1S,4R)-4-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl]cyclopent-2-  
en-1-amine;

(*S*) 1 {[4 (2,5-difluorophenyl) 2 phenyl 2,5 dihydro 1*H* pyrrol-1-yl]carbonyl}but-3-enylamine;

(*R*) 3 [4 (2,5-difluorophenyl) 2 phenyl 2,5 dihydro 1*H* pyrrol-1-yl] 3 oxo 1 phenylpropan-1-amine;

3 {[4 (2,5-difluorophenyl) 2 phenyl 2,5 dihydro 1*H* pyrrol-1-yl]carbonyl} 2 phenylpiperidine;

(*S*) 1 {[4 (2,5-difluorophenyl) 2 phenyl 2,5 dihydro 1*H* pyrrol-1-yl]carbonyl}but-3-enylamine;

(*S*) 3 [4 (2,5-difluorophenyl) 2 phenyl 2,5 dihydro 1*H* pyrrol-1-yl] 2 (methylamino)-3-oxopropan-1-ol;

(*R,S*) 5 {[4 (2,5-difluorophenyl) 2 phenyl 2,5 dihydro 1*H* pyrrol-1-yl]carbonyl}pyrrolidin-3-ol;

(*S*) 2 [4 (2,5-difluorophenyl) 2 phenyl 2,5 dihydro 1*H* pyrrol-1-yl] 2 oxo 1 (1,3-thiazol-4-ylmethyl)ethylamine;

(*R*) 1 {[4 (2,5-difluorophenyl) 2 phenyl 2,5 dihydro 1*H* pyrrol-1-yl]carbonyl}but-3-enylamine;

(*S*) 1 [4 (2,5-difluorophenyl) 2 phenyl 2,5 dihydro 1*H* pyrrol-1-yl] N,3 dimethyl-1-oxobutan-2-amine;

(*S*) 1 [4 (2,5-difluorophenyl) 2 phenyl 2,5 dihydro 1*H* pyrrol-1-yl] N,4 dimethyl-1-oxopentan-2-amine;

(*S*) 2 [4 (2,5-difluorophenyl) 2 phenyl 2,5 dihydro 1*H* pyrrol-1-yl] 1 [(1 methyl-1*H* imidazol-4-yl)methyl] 2 oxoethylamine;

4 (2,5-difluorophenyl) 1 (N-6 formyl L lysyl) 2 phenyl 2,5 dihydro 1*H* pyrrole;

(2S,3S)-1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,3-dimethyl-1-oxopentan-2-amine;

(1S)-1-(cyclohexylmethyl)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethylamine;

(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-(1H-indol-3-ylmethyl)-2-oxoethylamine;

(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-(isocyanoethyl)-2-oxoethylamine;

(1S)-1-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl]-3,3-dimethylbutylamine;

1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2,3-dimethyl-1-oxobutan-2-amine;

1-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl]cyclohexanamine;

1-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl]cyclopentanamine;

(1S)-3-(benzyloxy)-1-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl]propylamine;

1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2,3-dimethyl-1-oxobutan-2-amine;

1-[(4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl]cyclopent-3-en-1-amine;

(1S)-1-cyclopentyl-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethanamine;

4-(2,5-difluorophenyl)-1-(2-methylpropyl)-2-phenyl-2,5-dihydro-1H-pyrrole;

(1S)-2-[4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-cyclopropyl-2-exoethanamine;

(1S,2S)-1-[(4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl]-2-methylbutylamine;

(1S)-1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-but-3-enylamine;

(1S)-1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-but-3-ynylamine;

(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethanamine;

1-cyclopropyl-3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-exopropen-1-amine;

(1S,2S)-1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2-methylbutylamine;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-1-L-valyl-2,5-dihydro-1H-pyrrole;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-1-(2-methylalanyl)-2,5-dihydro-1H-pyrrole;

3-[1-[(2S)-2-amino-2-cyclopropylethanoyl]-4-(5-chloro-2-fluorophenyl)-2,5-dihydro-1H-pyrrol-2-yl]phenol;

~~4-(5-chloro-2-fluorophenyl)-2-(3-hydroxyphenyl)-1-L-isoleucyl-2,5-dihydro-1H-pyrrole;~~

(2S)-4-(2,5-Difluorophenyl)-N-methyl-2-phenyl-N-(piperidin-4-ylmethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(5-Chloro-2-fluorophenyl)-N-methyl-2-phenyl-N-[(3R)-pyrrolidin-3-yl]-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(5-chloro-2-fluorophenyl)-N-methyl-2-phenyl-N-piperidin-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(5-chloro-2-fluorophenyl)-N-methyl-2-phenyl-N-[(3S)-pyrrolidin-3-yl]-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-pyrrolidin-3-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-N-(1-allylpiperidin-4-yl)-4-(2,5-difluorophenyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-Difluorophenyl)-N-methyl-N-[(1-methylpiperidin-3-yl)methyl]-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-(pyridin-3-ylmethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-methyl-N-[(1-methyl-1H-pyrazol-4-yl)methyl]-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-[2-(dimethylamino)ethyl]-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-(2-pyridin-2-ylethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(5-chloro-2-fluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-piperidin-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-piperidin-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-1-[4-(2,5-difluorophenyl)-2-methyl-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-methyl-1-oxobutan-2-amine;

N-1-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-N-2,N-2-dimethylglycinamide;

N-1-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-N-2-methylglycinamide;

N-1-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)glycinamide;

N-1-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-methylalaninamide;

N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-pyrrolidin-1-ylacetamide;

2-azetidin-1-yl-N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)acetamide;

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-morpholin-4-ylacetamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-piperazin-1-ylacetamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-(4-methylpiperazin-1-yl)acetamide;~~

~~N-1-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-N-2-isopropylglycinamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)piperazine-1-carboxamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-N'-piperidin-4-ylurea;~~

~~4-amino-N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)piperidine-1-carboxamide;~~

~~N-(2-aminoethyl)-N'-(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)urea;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-N'-(3-morpholin-4-ylpropyl)urea;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-N'-(2-(dimethylamino)ethyl)urea;~~

~~2-azetidin-1-yl-N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)ethanesulfonamide~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-(isopropylamino)ethanesulfonamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-pyrrolidin-1-ylethanesulfonamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-morpholin-4-ylethanesulfonamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-piperazin-1-ylethanesulfonamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-(4-methylpiperazin-1-yl)ethanesulfonamide;~~

~~N-(tert-butyl)-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]acetamide;~~

~~2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-isopropylacetamide;~~

~~(2S)-1-(2-azetidin-1-yl-2-oxoethyl)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~(2S)-4-(2,5-difluorophenyl)-1-(2-oxo-2-pyrrolidin-1-ylethyl)-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~4-{[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]acetyl}morpholine;~~

~~1-{[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]acetyl}piperazine;~~

~~1-{[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]acetyl}-4-methylpiperazine;~~

~~2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-isopropylbutanamide;~~

4-{2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]butanoyl}morpholine;

2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-ethylacetamide;  
N-cyclobutyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]acetamide;

2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-ethylpropanamide;

N-cyclobutyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanamide;

2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-methylpropanamide;

2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-isopropylpropanamide;

N-(tert-butyl)-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanamide;

4-{2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanoyl}morpholine;

(3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-ethyl-2,2-dimethyl-4-oxobutanamide;

(3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2,2-dimethyl-4-oxo-N-piperidin-4-ylbutanamide;

(3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2,2-dimethyl-4-oxobutanoic acid;

(3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,N,2,2-tetramethyl-4-oxobutanamide;

(1S)-1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2,2-dimethyl-3-oxo-3-piperazin-1-ylpropylamine;

(3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-isopropyl-2,2-dimethyl-4-oxobutanamide;

(3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,2,2-trimethyl-4-oxobutanamide;

(3R)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,N,2,2-tetramethyl-4-oxobutanamide;

(3R)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2,2-dimethyl-4-oxobutanoic acid;

(1R)-1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2,2-dimethyl-3-oxo-3-piperazin-1-ylpropylamine

(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl-4-methylpiperazine-1-carboxylate;

(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl-1-methylpiperidin-4-ylcarbamate;

(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethylmethyl(1-methylpiperidin-4-yl)carbamate;

(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-exoethyl-4-dimethylamino)piperidine-1-carboxylate;

(2S)-N,N-dimethyl-4-(4-methylpyridin-3-yl)-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-{{[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]acetyl}morpholin-4-ium;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-*N*-methyl-*N*-piperidin-4-yl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

2-[{[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1*H*-pyrrol-1-yl]carbonyl}(methyl)amino]ethyl-4-methylpiperazine-1-carboxylate;

3-{4-(2,5-difluorophenyl)-1-[(4-methylpiperazin-1-yl)carbonyl]-2,5-dihydro-1*H*-pyrrol-2-yl}phenol;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-*N*-methyl-*N*-[(1-methyl-1*H*-pyrazol-4-yl)methyl]-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

4-(2,5-difluoro-phenyl)-2,5-dihydro-2-(3-hydroxyphenyl)-*N*-methyl-*N*-[2-(4-methyl-1-piperazinyl)-2-oxoethyl]-1*H*-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-*N*-methyl-*N*-[(5-oxo-4,5-dihydro-1*H*-1,2,4-triazol-3-yl)methyl]-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-*N*-{[5-(methoxymethyl)-1*H*-pyrazol-3-yl]methyl}-*N*-methyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-*N*-methyl-*N*-(1,3-thiazol-4-ylmethyl)-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-*N*-methyl-*N*-(1,3-thiazol-2-ylmethyl)-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-*N*-methyl-*N*-[2-(1*H*-1,2,4-triazol-1-yl)ethyl]-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-*N*-methyl-*N*-[2-(1*H*-pyrazol-1-yl)ethyl]-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

4-[*(2S)*-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]butan-2-one;

4-[*(2S)*-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-3-methylbutan-2-one;

3-[*(2S)*-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-*N*-methylpropanamide;

3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-*N,N*-dimethylpropanamide;  
3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-*N,N*,2-trimethylpropanamide;  
4-{3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]propanoyl}morpholine;  
1-{3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]propanoyl}-4-(methylsulfonyl)piperazine;  
1-{3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]propanoyl}piperidin-4-ol; and  
methyl 3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]propanoate.

16. (original) A pharmaceutical composition that is comprised of a compound in accordance with Claim 1 and a pharmaceutically acceptable carrier.

17. – 42. (Cancelled)